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R-51-5-2-15

**FINAL  
SCREENING SITE INSPECTION**

**LINFIELD INDUSTRIAL PARK**

**EPA WORK ASSIGNMENT NO. 37-34-3JZZ**

**PROJECT NO. 3263-09**

**EPA DSN PA-2898**

**FACILITY ID NO. PAD987332814**

**ARCS III PROGRAM**

**EPA CONTRACT NO. 68-W8-0037**

**SEPTEMBER 1992**



**HALLIBURTON NUS**  
*Environmental Corporation*

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ARCS III PROGRAM  
EPA CONTRACT NO. 68-W8-0037**

**FOR THE  
UNITED STATES ENVIRONMENTAL PROTECTION AGENCY**

**SEPTEMBER 9, 1992**

**SUBMITTED BY**

**(b) (4)**

**SITE MANAGER**

**REVIEWED BY**

**(b) (4)**

**PROJECT MANAGER**

**APPROVED BY**

**(b) (4)**

**PROGRAM MANAGER, ARCS III**

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## SECTION 1.0

## **1.0 INTRODUCTION**

### **1.1 AUTHORIZATION**

HALLIBURTON NUS Environmental Corporation performed this work under EPA Contract No. 68-W8-0037. This specific report was prepared in accordance with ARCS III Work Assignment No. 37-34-3JZZ for the Linfield Industrial Park Site, located in Linfield, Montgomery County, Pennsylvania.

### **1.2 SCOPE OF WORK**

HALLIBURTON NUS Environmental Corporation (formerly NUS Corporation) ARCS III was tasked to conduct a site inspection of the subject site.

### **1.3 SUMMARY**

The Linfield Industrial Park Site is located south of Linfield - Trappe Road, east of the Schuylkill River bridge, in Montgomery County, Pennsylvania. The residential town of Linfield is adjacent to and northeast of the site; state game lands are east of and adjacent to the site.

The abandoned 125-acre industrial park is situated on mostly flat land on the eastern banks of the Schuylkill River. The western and southern portions of the site have a five- to 10-percent slope toward the Schuylkill River. Access to the site is restricted by a fence and a main gate; however, the fence does not completely surround the site. The site is composed of approximately six separate divisions: the main gate and parking area, the distillery area, the building no. 2 area, the warehouse area, the concrete pads area, and the former tank farm area.

The site has been inactive since 1986. The current owner, 888 Warehousing, Incorporated, is attempting to find a lessee or buyer for the property.

From an unknown date until 1986, building no. 2 was used for the repackaging and distribution of ethylene glycol (antifreeze) and ammonia-based cleaners. These materials were transported to the facility by rail car and tanker trucks, transferred to smaller containers, and distributed off site for sale. Evidence at the site indicates that one-gallon plastic containers may have been manufactured on site for the distribution of the antifreeze and cleaner products.

Beginning in approximately 1945, the facility operated as a United-States-Government-bonded whiskey warehousing operation. The storage and distribution of whiskey comprised the main operation. Above-ground storage tanks provided the initial storage. Whiskey was then transferred to wooden kegs for aging in the warehouses. After the aging process, the whiskey was bottled and labeled at the site. It is not known how long this operation existed at the site.

For an undetermined period (probably beginning in the early 1800s), the facility was operated as a distillery of whiskey products. Storage, aging, and distribution of the whiskey products occurred at the site.

Inspections conducted by the Pennsylvania Department of Environmental Resources (PA DER) in 1972, 1973, 1975, 1983, 1984, and 1985 revealed unauthorized discharges of industrial wastewater to the Schuylkill River. The industrial wastewater consisted of deionizer wastewater, septic tank seepage, trash-compactor oils, bottle-making room oils, air compressor pit waste oil, and non-contact cooling compressor water. Additionally, the deionizer wastewater backwash and regeneration wastes were discharged after neutralization into a lime pit. From an unknown date until 1987, the facility was issued two NPDES permits for an outfall to the Schuylkill River. The NPDES discharge point was analyzed by PA DER on March 7, 1985 and revealed elevated levels of biochemical oxygen demand (BOD).

Groundwater and surface water are the sources of potable water for individuals residing in the four-mile-radius area surrounding the Linfield Industrial Park Site. The Citizens' Utilities Home Water Company (CUHWC) supplies water to a total population of 12,506 people in the study area. The water is obtained from an intake in the Schuylkill River (b) (9)

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The Pottstown Municipal Water Authority (PMWA)

supplies water to approximately 9,000 people in the study area. The water is obtained from an intake located upstream in the Schuylkill River, outside the study area. Approximately 19,831 persons are assumed to obtain their water from (b) (9)

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The Schuylkill River is adjacent to the southern and western boundaries of the site. PA DER lists the Schuylkill River as protected for the maintenance and/or propagation of fish species indigenous to warm water and for the passage, maintenance, and propagation of migratory fishes.

Adjacent to and downstream from the site are palustrine, forested, broad-leaved, deciduous, temporarily non-tidal flooded wetlands with three linear miles of frontage.

On February 19, 1991, HALLIBURTON NUS Field Investigation Team 3 (FIT 3) conducted a preliminary assessment of the Linfield Industrial Park. During the inspection, HALLIBURTON NUS FIT 3 observed several areas of concern: 21 above-ground storage tanks, two underground storage tanks that contained unknown oils, 75 to 100 fifty-five-gallon drums (many drums were empty, although at least 25 drums were not empty; 1 tar-like spill was observed), two power transformers, one gas pump, five tanker trucks (at least one truck was still hooked to the building via a discharge line), and 20 abandoned buildings. Based on these observations, EPA Emergency Response was contacted.

EPA Emergency Response performed an emergency assessment of the site on March 14, and 27, 1991. Before the assessment, the site owner informed the on-scene coordinator (OSC) that many of the drums observed by the FIT had been moved into an on-site warehouse. The large on-site holding tanks were found to be empty. Numerous drums and containers were found in the various on-site buildings. Four drums were sampled, and a soil sample was taken near a transformer. The OSC planned to work with the site owner to address safety concerns posed by the unrestricted site access and the materials haphazardly stored on site.

HALLIBURTON NUS ARCS III conducted a screening site inspection on January 30, 1992. Activities included sampling on-site surface soils and surface water and off-site groundwater and surface water. The results of the samplings, as shown in section 7.0 and discussed in section 8.0, revealed elevated levels of organic and inorganic contaminants including fluoranthene (72,000 ppb), benzo(a)pyrene (30,000 ppb), polychlorinated biphenyls (PCBs) (300,000 ppb), and lead (4,810 ppm) in on-site soil and elevated levels of inorganic contaminants including lead (59.90 ppb) and cyanide (11.70 ppb) in off-site surface water.

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## SECTION 2.0

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## 2.0 THE SITE

### 2.1 LOCATION

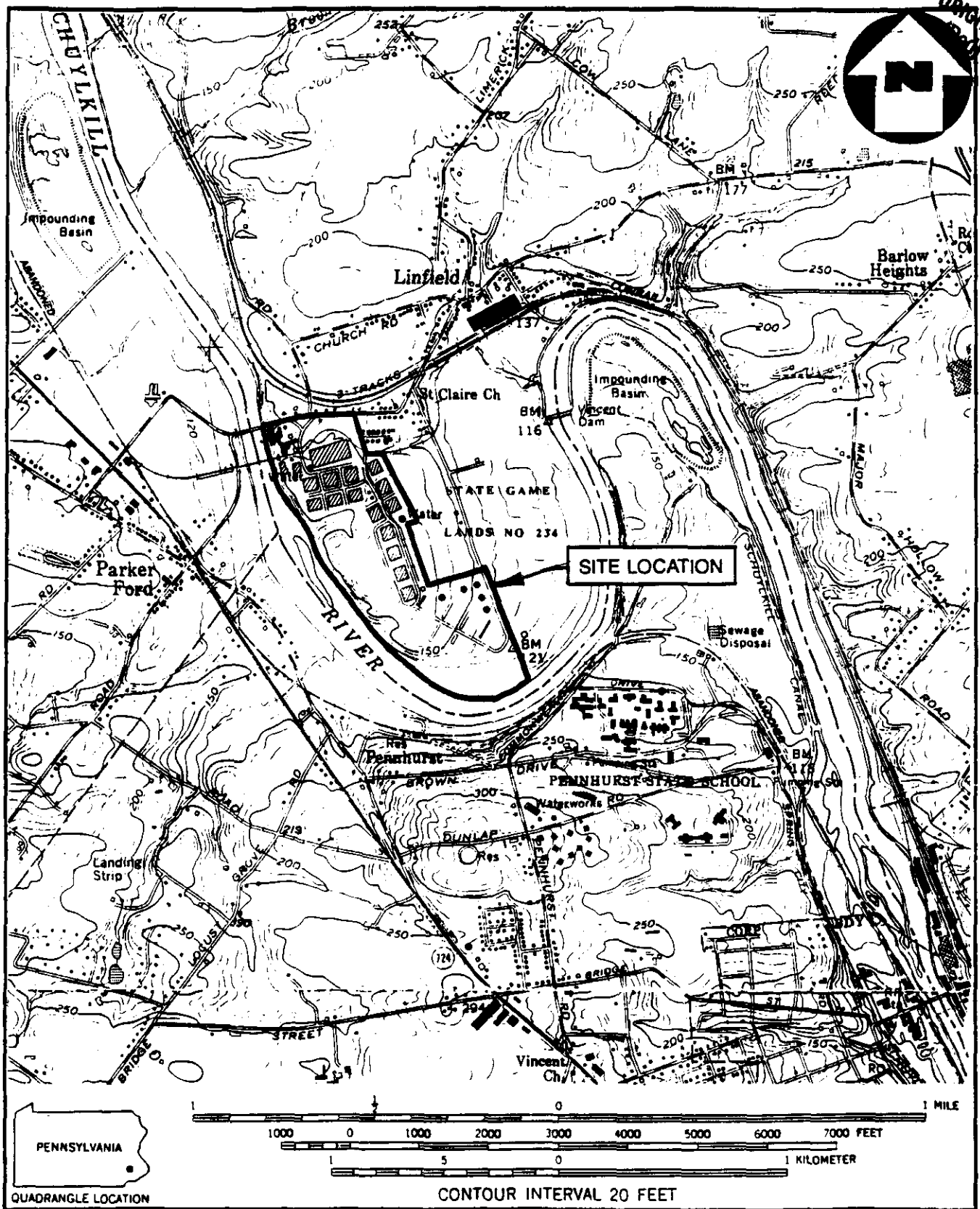
The Linfield Industrial Park Site is located immediately south of the Linfield - Trappe Road and east of the Schuylkill River bridge in Montgomery County, Pennsylvania (see figure 2.1, page 2-2). The site, which is south of Linfield, Pennsylvania, can be found on the United States Geological Survey (U.S.G.S.) Phoenixville, Pennsylvania 7.5 minute series topographic map at the coordinates 40° 12' 08" north latitude and 75° 34' 38" west longitude. As measured from the southwestern corner of the Phoenixville, Pennsylvania topographic map, the site is approximately 6-3/4 inches east and 14 inches north.<sup>1,2,3</sup>

### 2.2 SITE LAYOUT

The 125-acre industrial park site is located immediately off Linfield - Trappe Road, south of Linfield, Pennsylvania. The site is bordered to the south and west by the Schuylkill River, to the east by State Game Land No. 234, and to the north and northeast by the residential town of Linfield. ConRail Corporation railroad tracks are located in the town of Linfield, north of the site; abandoned tracks are throughout the site (see figure 2.2, page 2-3).<sup>2</sup>

The site can be divided into six divisions: the fenced parking lot, the distillery area (distillery building, office/storehouse, abandoned house, leachate field, and other buildings), building no. 2 or the ethylene glycol area, the warehouses, the concrete pads, and the former tank farm.<sup>2</sup>

Located on the northeastern portion of the site is a paved access road that leads from Linfield - Trappe Road to a 200- by 75-foot parking lot on site. A guardhouse was adjacent to the main gate. A chain-link fence separated the parking lot from the facility. An abandoned tanker truck was northwest of the guardhouse at the time of the preliminary assessment.<sup>2</sup> The guardhouse was not occupied at the time of the site inspection.<sup>4</sup>



SOURCE: (7.5 MINUTE SERIES) U.S.G.S. PHOENIXVILLE, PA. QUAD.

**SITE LOCATION MAP**  
**LINFIELD INDUSTRIAL PARK, LINFIELD, PA.**

SCALE 1: 24000

FIGURE 2.1







Located on the northwestern portion of the site was the two-story former distillery building. At the time of the preliminary assessment, the distillery and the area surrounding it consisted of eight whiskey vats, oil and grease drums, five above-ground storage tanks, an oven and smoke stack with a coal pile, a transformer, and abandoned trucks. The distillery building housed the equipment used in the whiskey manufacturing process, including approximately eight empty (11,700-gallon capacity each) wooden fermentation vats. Adjacent to the distillery building on the northern side were five above-ground metal cylinder storage tanks. Three of these tanks hold approximately 5,284 gallons each; the two remaining tanks hold approximately 939 gallons each. The contents of these tanks are not known. Adjacent to and east of these tanks were an electrical transformer unit, and an area of stained soils was noted. Several 55-gallon drums of oil and grease were observed in this area during the preliminary assessment. On the southern side of the distillery building was a coal pile that was approximately 10 by 15 by six feet in size (150 square feet of ground surface area was covered). A stoker oven, east of the coal pile, had piping and a smoke stack attached to the southern side of the building.<sup>2</sup>

North of the distillery was an abandoned house. Adjacent to the eastern side of the house was an inactive well, and south of the house was a gas pump.<sup>2</sup>

An office/storehouse building was located northwest of the distillery. Pipes led to the Schuylkill River from the western side of the building. Two underground storage tanks, approximately 10,000 gallons each, containing oil-like liquids were located east of the building; the underground storage tanks could be accessed through manhole covers. At the time of the preliminary assessment, several 55-gallon drums were located near the southern side of the building; at least two drums were full and marked "Used Oil." At the bottom of the hill, adjacent to the Schuylkill River and west of the building, was a concrete underground open pit that appeared to be a surface water intake. A pump and pipeline system with a broken screen on top were observed.<sup>2</sup>

South of the office/storehouse and west of the distillery building was a leachate field with vent stacks. A 12- by 12-foot building was located in this area. A small generator with indications of stained soils was observed in this area. Ten 55-gallon drums of oil and varnish were located inside the building. West of the leachate field was the former NPDES outfall. South of the outfall, abandoned railroad tracks crossed over the river.<sup>2</sup>

South of the distillery building were two buildings and several abandoned trucks. At the time of the preliminary assessment, the eastern building contained personnel lockers and pumping equipment. The contents of the western building are not known. A large above-ground tank, 30 by 50 feet in size, with an approximately 264,231-gallon capacity, was located adjacent to the buildings. The trucks in this area were a New York Fire Department truck and several trailer trucks.<sup>2</sup>

In the north-central portion of the site was building no. 2 or the ethylene glycol area. This building was reportedly used primarily for ethylene glycol bottling and distribution. East of the building were 10 fenced, black, above-ground metal cylinder storage tanks, six by 40 feet in size, with an 8,455-gallon capacity. The contents of these tanks are not known. One white above-ground metal cylinder storage tank, five by 30 feet in size, with an approximately 4,400-gallon capacity, was located on the northeastern side of the building. This white tank contained plastic white pellets. These white pellets were also on the ground around the tank. One brushed aluminum above-ground metal cylinder storage tank, five by 10 feet in size, with a 1,467-gallon capacity, was located here also. A white above-ground propane storage tank was located 200 feet north of the building. A loading dock was on the northern side of building no. 2. Four abandoned tanker trucks were observed adjacent to the building. One truck on the southern side of the building was connected to the building via its discharge pipe. The contents of the tankers are not known. West of building no. 2 was a water tower.<sup>2</sup>

Located centrally on the site and south of building no. 2 was the warehouse area. The 14 warehouse buildings (three-story and windowless) were formerly used for the storage of whiskey barrels. Two above-ground aluminum cylinder storage tanks, 40 by 50 feet in size, with an approximately 469,244-gallon capacity, and one above-ground metal cylinder tank, eight by 15 feet in size, with a 3,637-gallon capacity, were located centrally among these warehouses. The contents of these tanks are not known. A water tower was located in this area. Pipelines insulated in asbestos-like material (some pipelines had collapsed) connected the warehouses.<sup>2</sup>

During the site inspection, a trash dump measuring approximately 30 by 75 feet and consisting mostly of empty plastic antifreeze containers was observed between the southernmost warehouse and the Schuylkill River. Adjacent to this area were approximately 60 upright empty and rusted drums. No markings were visible on the drums.<sup>4</sup>

A drainage swale that led to the Schuylkill River was in a low-lying area west of the warehouses in the middle of the site. The swale originated at the outfall of a large concrete pipe. Several smaller pipe outfalls and three concrete-lidded manholes were also in this area. In an open field east of this area was an approximately 20- by 100-foot area of stressed vegetation.<sup>4</sup>

Located on the southwestern part of the site were eight concrete pads. These 30- by 250-foot concrete roads extended to the east and west from a main north-south-trending on-site road. Demolished wooden buildings were observed at each pad area. At least twenty-two 55-gallon drums contained an unknown liquid at the time of the preliminary assessment. Four drums were near the northwestern pad area, and eight drums were located near the southeastern pad area. One 55-gallon drum contained an oil-like residue, which had spilled and stained the ground. Several dump sites on these pads consisted of antifreeze containers, piles of antifreeze jugs and melted plastics, roofing materials, and miscellaneous wastes. Throughout the area were whiskey bottles and wooden barrels. A 12- by 12-foot building was observed west of the main road and between two concrete pads. Five 5-gallon containers marked "Herbicide" were observed in this building during the preliminary assessment. The contents of the containers are not known. Southeast of the concrete pad areas was an abandoned house. At this location, a small electrical transformer was observed.<sup>2</sup>

During the site inspection, an approximately 100- by 100-foot area of stressed vegetation was observed northwest of the abandoned house. This area was quantitatively identical to the stressed vegetation west of the middle warehouses.<sup>4</sup>

Located northeast of the concrete pad areas was a former tank farm. Five former tanks areas, which were individually fenced and bermed in concrete, were located in this area of the site. The tanks were not present during the preliminary assessment; however, the concrete containment areas still existed.<sup>2</sup>

### **2.3 OWNERSHIP HISTORY**

The Linfield Industrial Park Site has been owned by 888 Warehousing, Incorporated, of Brooklyn, New York, since 1986. The park is closed and is not utilized for any purpose.<sup>2,5</sup>

Before 1986, the site was owned by Publicker Industries, Incorporated, of Philadelphia, Pennsylvania. The site representative, Bernard Shafran, stated that Mr. Publicker had owned the site since at least sometime after World War II. During the ownership by Publicker, the site was operated under different names: Continental Distilling Corporation, Publicker Packaging Services, Publicker Industries, and Linfield Industrial Park. There is no record to indicate who owned the park before Mr. Publicker.<sup>6,7,8</sup>

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Site Name: Linfield Industrial Park  
Project No.: 3263-09

ConRail Consolidated Corporation, of Philadelphia, Pennsylvania, apparently owns the rail lines throughout the site. According to the HALLIBURTON NUS FIT 3 Limerick Township site discovery, the tax office of Limerick, Pennsylvania shows the property tax as being paid by ConRail.<sup>3</sup>

## **2.4 SITE USE HISTORY**

The site has been inactive since June 1986. The current owner, 888 Warehousing, Incorporated, has no immediate plans for the site unless a lessee or buyer can be found.<sup>2,5</sup>

For an undetermined time until 1986, building no. 2 was used to repackage and distribute ethylene glycol (antifreeze) and ammonia-based cleaners. These materials were transported to the facility by rail car and tanker trucks, transferred to smaller containers, and distributed off site for sale. Evidence at the site indicates that one-gallon plastic containers for antifreeze and cleaners may have been manufactured on site.<sup>2,3,9</sup>

For an undetermined time beginning in approximately 1945, the facility was operated as a United-States-Government-bonded whiskey warehousing operation. The storage and distribution of whiskey were the main operations. Above-ground storage tanks were the initial means of storage. Whiskey was then transferred to wooden kegs for aging in the warehouses. HALLIBURTON NUS FIT 3 observed that many wooden kegs in various stages of construction and wooden keg pieces were stockpiled, indicating that these wooden kegs were built on site. After the aging process, the whiskey was bottled and labeled at the site. The bottled whiskey product was distributed by rail cars and trucks.<sup>2</sup>

The facility was once operated as a whiskey distillery. Although the dates of this operation are not known, the distillery building appears to have been constructed in the early 1800s. Large vats were present in the distillery building. Storage, aging, and distribution of the whiskey products were the main operations.<sup>2</sup>

The site use before the distillery was constructed is not known.

## 2.5 PERMIT AND REGULATORY ACTION HISTORY

Permit and regulatory action relevant to the subject site is discussed chronologically below:

On September 5, 1972, a violation of the Clean Streams Law was documented by PA DER after an inspection on August 15, 1972 revealed several discharges from the company to the Schuylkill River without a permit. These discharges included deionized wastewater, cooling water, sink wastewater, water-softener waste, and septic-tank seepage.<sup>8</sup> According to PA DER correspondence dated January 22, 1974, on January 11, 1973, PA DER sent information to Continental Distilling Corporation (the operating name of the site at this time) detailing the information needed to apply for the industrial wastewaters. No record of results was found. On October 31, 1973, an inspection by PA DER revealed that the facility was still in violation of the Clean Streams Law because of unauthorized discharges. On January 22, 1974, PA DER acknowledged that no information had been received to date from Continental Distilling Corporation regarding permit application.<sup>10</sup>

On February 7, 1974, an administrative conference was held between PA DER and Continental Distilling Corporation. During the conference, Continental Distilling submitted the NPDES permit project status cards. On February 8, 1974, PA DER determined that a payment of \$250 to the Clean Water Fund by Continental Distilling Corporation would be considered as final settlement of all claims for the violations of the Clean Streams Law noted during the inspection conducted on February 27, 1973.<sup>11</sup>

On February 10, 1975, PA DER requested updated NPDES permit project schedule cards and the completion of a permit application.<sup>12</sup> According to correspondence dated July 10, 1975, on February 26, 1975, PA DER had acknowledged the receipt of NPDES permit schedule cards from Continental Distilling Corporation. The NPDES permit application was not submitted.<sup>13</sup>

On April 21, 1975, PA DER conducted an inspection of the Continental Distilling Corporation facility; the inspection revealed that the facility's septic system was overflowing. Samples were obtained during the inspection. One discharge from the site had a BOD of more than 80 ppm, and another discharge from the site had a BOD of more than 300 ppm and an ethyl glycol content more than 280 ppm.<sup>13</sup> The BOD results indicate the amount of dissolved oxygen used up by the sample when incubated in darkness at 20°C for five days.<sup>14</sup> It is not documented who took the samples or where the samples were analyzed. PA DER considered both discharges from the site to be violations of the Clean Streams Law because they were unpermitted discharges.<sup>13</sup>

On July 10, 1975, PA DER requested letter notification from Continental Distilling Corporation of the status of its NPDES permit application, the treatment or elimination of contaminated waste discharges, and the facility's plan to correct the septic system problem.<sup>13</sup>

On March 19, 1976, an administrative conference was held between PA DER and Continental Distilling Corporation. The following topics were covered: tile field overflows, unpermitted industrial waste discharges, NPDES application, NPDES permit project status schedule cards, and the Pollution Incident Prevention Plan (PIPP). PA DER required that the old septic tank connected with the old tile field be pumped out as soon as the new tile field was in operation. PA DER requested that updated NPDES permit project status schedule cards and an NPDES permit application be submitted. No specifics were stated regarding the PIPP.<sup>15</sup>

On April 6, 1976, NPDES project status schedule cards were accepted by PA DER.<sup>16</sup> According to correspondence dated July 13, 1979, on May 18, 1979, PA DER requested a resubmission of the industrial waste permit application for reissue of NPDES Permit No. PA0013293.<sup>17</sup>

On July 13, 1979, PA DER required Continental Distilling Corporation to submit a Part II (or better) permit for the discharges to the groundwater via the limestone pit adjacent to warehouse no. 10 and recommended that a treatment facility be used for neutralization and suspended solids removal.<sup>17</sup>

On January 11, 1983, PA DER notified Continental Distilling Corporation that the expiration date of NPDES Permit No. PA0013293 was July 12, 1983.<sup>18</sup>

On April 27, 1983, PA DER conducted an inspection that revealed the discharge of non-contact compressor cooling water to a storm sewer tributary to the Schuylkill River. On May 5, 1983, PA DER advised Publicker Industries that such unauthorized surface water discharges were occurring. The Part I NPDES application was enclosed with this letter.<sup>19</sup>

On January 22, 1985, PA DER issued a letter to Publicker Industries regarding a December 6, 1984 inspection that had been conducted by PA DER in response to citizens' complaints that eye and nasal irritation occurred to persons walking near the cooling water discharge. The inspection revealed that the company was processing "CLEAN JOB," an ammonia-based cleaner. Samples obtained during the visit revealed a BOD more than 405 ppm and phenol concentrations of 5.0 ppb.<sup>20</sup>

On May 8, 1985, PA DER confirmed the results of an inspection that had been conducted by PA DER on March 7, 1985. Samples obtained during the inspection revealed that the non-contact cooling water discharged to outfall no. 001 contained a BOD concentration of 11.5 ppm, which violated the terms of NPDES Permit No. PA0013293. The source of contamination was not determined.<sup>21</sup>

On August 13, 1985, PA DER confirmed the results of an inspection that had been conducted by PA DER on July 17, 1985. The inspection results revealed the following sources of potential pollution at the site that could impact the outfall no. 001 discharge to the Schuylkill River: a floor drain in the blending room leading to outfall no. 001, oil tankers with stained soils into the storm sewer, oils in air compressor pit discharge into outfall no. 001, trash compactor oils near the storm sewer inlet, and oil leaks in the bottle room that were discharged to the ground. These conditions were determined by PA DER to be violations of the Clean Streams Law. No Notices of Violation or penalties were filed, however.<sup>22</sup>

On May 11, 1987, PA DER requested a renewal application for Publicker Packaging Company's NPDES Permit No. PA0013293; the current permit would expire on December 23, 1988.<sup>23</sup>

On February 19, 1991, HALLIBURTON NUS FIT 3 conducted a preliminary assessment inspection of the Linfield Industrial Park Site.<sup>2</sup>

EPA Emergency Response performed an emergency assessment of the site on March 14 and 27, 1991. Many of the drums observed by the FIT had been moved into an on-site warehouse by the site owner. During the first EPA Emergency Response site visit, during which on-site buildings were not accessed, the following were noted on site: five drums containing material, four truck trailers, three tankers, and an abandoned barn containing transformer carcasses. Stained soil near a transformer field tested positive for PCB contamination greater than 50 ppm. During the second site visit, on-site buildings were investigated. Numerous drums and containers were found inside the various buildings, including containers of sulfuric acid, drums and containers marked "Flammable," containers of liquid caustic materials, thousands of small containers filled with antifreeze, drums of methanol, and a drum marked "Toluene." The buildings were secure and locked. Five samples were taken during the assessment: four were drawn from drums and one was taken from the soil near the transformer. The results of the sample analyses are not known. The large on-site holding tanks were found to be empty. No removal or enforcement action was conducted.<sup>24,25,26</sup>



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## **2.6 REMEDIAL ACTION TO DATE**

On September 9, 1985, Publicker Packaging Services contacted PA DER regarding potential pollution sources revealed in PA DER's August 1985 inspection of potential impacts on NPDES outfall no. 001 to the Schuylkill River. The following comments were offered by Publicker Packaging Services:<sup>7,27</sup>

- The floor drains in the blending area had been sealed off from the blending pit since early 1985. The drains empty into a sump from which material is pumped into a holding tank for disposal. There is a by-pass valve in the line to the holding tank, which, if opened, would allow material to flow into the blending pit. This branch was sealed off to prevent any accidental discharges.
- The ground area at the bottle-making room was immediately dug out. The soil had been disposed through the Boyertown Landfill facility. The quantity of removed soil is not known.

On July 14, 1975, Continental Distilling Corporation contacted PA DER regarding ethyl alcohol found in one of the discharges. The following comment was offered by Continental Distilling Corporation:<sup>28</sup>

- A contractor was proposed to be hired for the removal of the existing tile field and the excavation of 150 square feet of soil and for the installation of a new stone base and tile, cover, and fill. It is not known if these activities were completed. The quantity of contaminated soil removed is not known.

On December 23, 1985, Publicker Packaging Services contacted PA DER regarding corrective actions taken in response to a spill on December 14, 1985. The following comments were offered by Publicker Packaging Services:<sup>29</sup>

- All residual antifreeze had been vacuumed or picked up with absorbent material. Eldredge Waste Management, of West Chester, Pennsylvania, had flushed and removed residual antifreeze from the drain system. A valve had been installed to prevent any liquids from entering the pipeline. The quantity of contaminated material removed is not known.

No other remediation has taken place at the site.

ORIGINAL  
2000

### SECTION 3.0

### 3.0 ENVIRONMENTAL SETTING

#### 3.1 WATER SUPPLY

Residents within the study area rely on surface water and groundwater sources for their potable water supplies. CUHWC serves residents in and around the boroughs of Spring City and Royersford within the study area and relies on surface water and groundwater sources of supply. PMWA serves residents in the northwestern part of the study area and relies upon a surface water source of supply. PBWS serves a very small number of residents in the extreme southeastern part of the study area and relies upon a surface water source of supply. The remaining residents of the study area are assumed to maintain private wells or springs for their drinking water supply. 1,30,31,32

CUHWC supplies water within the boroughs of Spring City and Royersford and to small areas of neighboring East Coventry, East Vincent, Limerick, and Upper Providence Townships. Water for this system is obtained from an intake on the Schuylkill River and from three wells. Some water is also purchased from PBWS. The Schuylkill River intake is located directly across the river from the site

(b) (9)

(b) (9)

(b) (9)

Water source withdrawal and purchase data for

report year 1989 are shown below. 1,30,31,32,33,34,35,36, 37

CUHWC Source	Withdrawal/ Purchase (gallons per day)	Number of Days	1989 Total	Percent of System (1989 Total)
Schuylkill River	817,000	248	202,616,000	38.2
(b) (9)	210,000	306	64,260,000	12.1
(b) (9)	302,000	121	36,542,000	6.9
(b) (9)	612,000	350	214,200,000	40.4
Purchase PBWS	35,100	365	12,811,500	2.4
<b>TOTAL</b>	<b>1,453,000</b>	<b>365</b>	<b>530,429,500</b>	<b>100.00</b>

The CUHWC system is fully integrated and supplies water to a total population of 12,506 persons 1,30,31,32,33,34,35,36,37

PMWA supplies water to the borough of Pottstown and to portions of neighboring North Coventry and Lower Pottsgrove Townships within the study area. PMWA draws 100 percent of its water supply from an intake on the Schuylkill River located (b) (9) PMWA does not purchase water from or sell water to other public water suppliers. A total population of 31,500 persons is supplied with water by PMWA. 1.30.31.32.37,38

The Phoenixville Borough Water System (PBWS) supplies water to the borough of Phoenixville, the community of Mont Clare, and small areas of East Pikeland and Schuylkill Townships. PBWS draws 100 percent of its water supply from an intake on the Schuylkill River (b) (9) (b) (9) PBWS supplies water to a total population of 20,000 persons and sells water to CUHWC. 1.32.37,39,40

The Philadelphia Suburban Water Company (PSWC) maintains a surface water intake on the Schuylkill River (b) (9) PSWC maintains an integrated system supplying water to 222,660 residential connections in areas of Chester, Delaware, and Montgomery Counties. Water for the integrated system is drawn from 37 active wells, five surface water intakes, and a groundwater reservoir located in a former quarry. Water is also purchased from the Pennsylvania - American Water Company (PAWC) - Norristown District and the West Chester Area Municipal Authority. Water source withdrawals and purchase data for report year 1990 are summarized below. 1.41.42.43,44,45, 46

PSWC Source	Withdrawal/ Purchase (gallons per day)	Number of Days	Percent of Total Purchases and Withdrawals
37 wells combined	13,670,000	365	14.8
Upper Merion Reservoir	6,160,000	365	6.7
Crum Creek	20,000,000	365	21.6
Pickering Creek	5,420,000	365	5.9
Perkiomen Creek	19,500,000	365	21.1
Neshaminy Creek	9,590,000	365	10.4
Schuylkill River	11,400,000	365	12.3
Purchases	6,660,000	365	7.2
<b>TOTAL</b>	<b>92,400,000</b>	<b>365</b>	<b>100.00</b>

PSWC also sells water to the Malvern Borough Water Department, PAWC Norristown District, the North Wales Water Authority, and the Hatboro Borough Authority. 1.41.42.43,44,45,46

The majority of the private domestic wells in the area produce water from the (b) (9) Formation. The remaining wells produce from the (b) (9) (b) (9) Wells in the (b) (9) Formation of Limerick Township have median depths of (b) (9) and a median yield of 15 gallons per minute (gpm). (b) (9) (b) (9) (b) (9) 1,30,32,36,47

### 3.2 SURFACE WATERS

The Schuylkill River is adjacent to the southern and western boundaries of the site.<sup>1,2</sup> Surface drainage from the site is expected to travel southwestward and flow into the Schuylkill River.<sup>2</sup> The former NPDES outfall is in the northwestern part of the site. Violations of the Clean Water Act have been found during inspections of this outfall by PA DER since 1972.<sup>2,8,10,12, 13, 15, 16, 17, 22</sup>

PA DER lists the Schuylkill River as protected for the maintenance and/or propagation of fish species indigenous to a warm-water habitat and for the passage, maintenance, and propagation of migratory fishes.<sup>48</sup> The Schuylkill River has an average flow of 1,888 cubic feet per second (cfs).<sup>49</sup> The site is located within the 100-year flood-prone area of the Schuylkill River.<sup>50</sup>

Immediately downstream from the site are palustrine, forested, broad-leaved, deciduous, temporarily non-tidal flooded wetlands with approximately three linear miles of frontage.<sup>51</sup>

CUHWC utilizes an intake in the Schuylkill River. (b) (9)

PBWS maintains a surface water intake (b) (9)

PSWC maintains a surface water intake (b) (9)

### 3.3 HYDROGEOLOGY

The geologic and hydrogeologic conditions in the study area were researched as part of the site inspection. A preliminary literature review was conducted to determine surface and subsurface geologic conditions, soil character, and the status of groundwater transport and storage.

### **3.3.1 Geology**

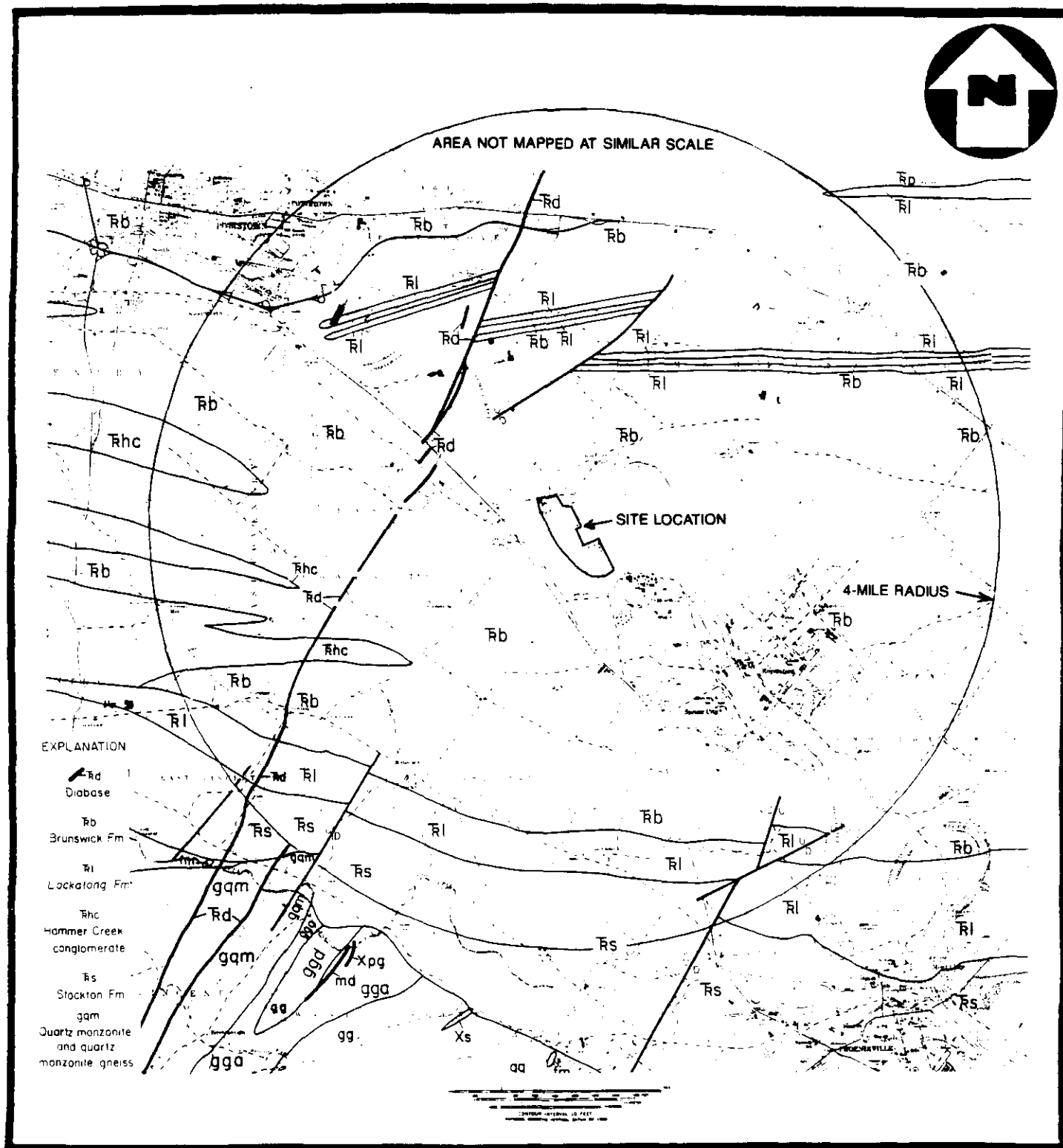
The site lies within the Triassic Lowlands Section of the Piedmont Physiographic Province. The area is predominantly an undulating plain with relief provided by low hills and ridges. These hills and ridges were formed by the differential weathering and erosion of the sandstones and shales that underlie most of the area. Steeper slopes may be formed along the sides of the Schuylkill River valley, which has been cut below the level of the surrounding countryside. The drainage pattern of the area is dendritic, and the entire study area is drained by the Schuylkill River and its tributaries.<sup>1,52,53</sup>

Virtually all of the study area is underlain by rocks of the Late Triassic age Newark Group, which includes the Lockatong, Hammer Creek, Brunswick, and Stockton Formations. These sedimentary rocks have been intruded by diabase dikes and sills, also of Triassic age. These rocks are part of a large Triassic basin that trends generally northeastward from Virginia to New York. The average dip of the beds within this basin is to the north or northwest at about 20 degrees. Within the vicinity of the site, the bedding dips mainly to the north at about 15 degrees. A very small portion of the southwesternmost part of the study area is underlain by quartz monzonite and quartz monzonite gneiss of the Piedmont Uplands Section of the Piedmont Physiographic Province.<sup>36,54,55,56</sup>

The rocks of the area are cut by many faults, most of which are relatively small. The nearest of these with respect to the site is located one or more miles to the north. This fault strikes to the northeast and is downthrown to the southeast, noticeably offsetting the dipping units of the Brunswick and Lockatong Formation shown on the geologic map (see figure 3.1, page 3-5). Joint systems are well developed in many of the beds of the Brunswick Formation. A very small set of joints strikes about north 30 degrees east. One or both of two less well-developed joint sets may be present at most locations. These sets strike about north 45 degrees west and north 75 degrees east. All joints are nearly vertical, with an average spacing of about six inches. Their orientation appears to be independent of the strike and dip of the beds.<sup>36,52,54</sup>

The formation directly underlying the site and most of the study area is the Triassic age Brunswick Formation. It consists mainly of reddish-brown shale, mudstone, and siltstone. A few very thin beds of green or brown shale are present in some places. Tough, thick-bedded, red argillite is interbedded with a dark gray argillite, typical of the Lockatong Formation, near the base of the Brunswick. Many of the shale beds are micaceous, causing them to split evenly along bedding planes. Calcite and quartz and occasionally barite and pyrite are present, partly filling joints within the Brunswick. The beds of the Brunswick Formation strike from west to east and have dips of approximately 14 to 15 degrees to the north in the vicinity of the site. The exact strike and dip of the beds underlying the site are not known. The Brunswick Formation attains a maximum thickness of 9,000 to 16,000 feet; however, the exact thickness beneath the site is unknown.<sup>36,52,54,57</sup>

ORIGINAL  
(Red)



SOURCE: Pennsylvania Geologic Survey. Atlas of Preliminary Geologic Quadrangle Maps of Pennsylvania. Map 61. 1981.

**GEOLOGIC MAP**  
**LINFIELD INDUSTRIAL PARK**  
**MONTGOMERY CO., PA.**

3-5

**FIGURE 3.1**



**HALLIBURTON NUS**  
*Environmental Corporation*

ORIGINAL  
FILE

Stratigraphically underlying and occurring within the Brunswick Formation is the Triassic Lockatong Formation. The nearest estimated subcrop or outcrop locations of the Lockatong Formation occur 1.5 miles or more north and 2.9 miles or more south of the site. The formation consists mainly of medium to dark gray argillite interbedded with thin beds of gray to black shale, siltstone, and marlstone. This formation is usually thick bedded or massive. It is made up largely of analcime, dolomite, feldspar, and clay with scattered pyrite. Calcite is common, especially filling joints, and quartz is a very minor constituent. The Lockatong has a maximum thickness of about 1,500 feet in this area near the Schuylkill River.<sup>36,52,54</sup>

Interfingering with the Brunswick Formation is the Hammer Creek Conglomerate. The nearest estimated subcrop or outcrop locations of this unit occur 1.8 or more miles southwest and west of the site. The Hammer Creek is equivalent in age to the Brunswick, and the division of the rocks into these two units is based mainly upon lateral changes in lithologic character. Rocks of the Hammer Creek Conglomerate are much more coarse than those of the Brunswick Formation. They consist of very coarse quartz conglomerate with abundant pebbles and cobbles of gray quartzite along with minor interbeds of coarse red sandstone. Based on outcrop patterns and an average dip of 20 degrees, the Hammer Creek Conglomerate is probably 500 feet thick or less within the study area.<sup>36,58,59</sup>

Stratigraphically underlying the Lockatong Formation about 3.2 miles south of the site is the Triassic Stockton Formation. The estimated subcrop or outcrop location occurs 3.2 miles south of the site. It is the oldest of the Triassic rocks found within the study area and is divided into three members. The lower arkose member is characterized by an abundance of coarse-grained arkosic sandstone and arkosic conglomerate. The middle arkose member is characterized by an abundance of fine- and medium-grained arkosic sandstone. The upper shale member consists mainly of shale and siltstone. The sandstones of the Stockton are yellow and brown to red, and the shales are red and soft. Sediments of the different textures are irregularly bedded, and beds commonly pinch out or grade laterally into beds of different texture and color. The Stockton Formation is about 2,300 feet thick in the Phoenixville area.<sup>36,55</sup>

Intruding the Brunswick, Lockatong, Hammer Creek, and Stockton Formations is a series of Triassic diabase sills and dikes. The nearest of these is located about 1.2 miles west-northwest of the site. The diabase intruded as dikes is black, dense, and very fine grained. It consists mainly of labradorite and augite. The dikes are typically five to 100 feet thick. The shales of the Brunswick Formation have been altered to a dark, tough hornfels close to the diabase intrusives. These alteration zones vary greatly in width but are usually between 40 and 100 feet in the vicinity of the smaller dikes.<sup>36,52,54</sup>



Underlying the Stockton Formation approximately 3.9 miles southwest of the site are quartz monzonite and quartz monzonite gneiss of probable Precambrian age. The unit is a medium- to fine-grained, medium to dark gray rock composed primarily of feldspar and quartz, with dark accessory and alteration minerals including pyroxene, amphibole, and biotite. Locally, the rocks show a foliated or gneissic texture. The thickness of this unit is not estimated.<sup>57,60</sup>

### **3.3.2 Soils**

The soils underlying the site are classified as Made land, Birdsboro, Penn, and Rowland silt loams, and Stony land. The distribution of these units is shown on figure 3.2 (page 3-8). The native soils have been disturbed and/or covered throughout large portions of the site by the construction of buildings, tanks, railroad spurs, roads, and other site facilities.<sup>2,61</sup>

Made land (MeB on figure 3.2) covers approximately 50 percent of the site in the northern and central-eastern portions. This unit is more specifically described as Made land, shale and sandstone materials, sloping. It results from the altering and mixing of soils that formed in material weathered from shale and sandstone. It is mainly nearly level and gently sloping but also includes some moderately sloping and steep areas. It is used for residential, industrial, commercial, and institutional development. Dusky-red to yellowish-brown shaly silt loam to channery sandy loam comprise much of the unit, and many areas consist entirely of pieces of shale. Bedrock may crop out in some places and may be as much as six feet deep in other places. The estimated permeability ranges from moderate to very slow (2.0 to less than 0.2 inch per hour). The soil reaction is medium acid to very strongly acid (pH, 6.0 to 4.5).<sup>61</sup>

Birdsboro silt loam soils cover approximately 20 percent of the site in the southeastern portion. They are present as the Birdsboro silt loam, zero to three percent slopes (BuA on figure 3.2) and as Birdsboro silt loam, three to eight percent slopes, moderately eroded (BnB2 on figure 3.2). They are deep, well-drained, reddish-brown soils that formed from stream sediments washed from uplands underlain by red shale and sandstone. In a typical Birdsboro soil profile, the surface layer is friable, dark, reddish-brown silt loam with a few river pebbles. The surface layer is about eight inches thick. The subsoil is friable, reddish-brown silty clay loam or clay loam 2-1/2 to three feet thick. The substratum is red or reddish-brown sandy loam about two feet thick and contains considerable gravel. The substratum is underlain by a layer of firm, dusky-red silt loam weathered from underlying rock. Soft, weathered dusky red shale bedrock is typically found at a depth of seven feet; however, the depth to bedrock ranges from four to 15 feet. Birdsboro soils have moderate permeability (0.63 to two inches per hour) and high available moisture capacity. The soil reaction is strongly acid to medium acid (pH, 5.1 to 6.0).<sup>61</sup>



SOILS MAP  
 LINFIELD INDUSTRIAL PARK  
 MONTGOMERY CO., PA.

Rowland silt loam, coal overwash (Ru on figure 3.2), covers approximately 10 percent of the site along the western and southern boundaries adjacent to the Schuylkill River. Rowland soils are deep, moderately well-drained to somewhat poorly drained, nearly level silt loams on floodplains. They formed from material washed from uplands underlain by red shale and sandstone. The Rowland silt loam, coal overwash unit has a surface layer of black silt loam that is one to three feet thick. This is underlain by a substratum consisting of several layers of dark reddish-gray silt loam. Shale or sandstone bedrock occurs at a depth of three to 12 feet. This soil has moderate permeability (0.63 to two inches per hour) and high available moisture capacity. The soil reaction is very strongly acid to medium acid (pH, 4.5 to 6.0).<sup>61</sup>

Penn silt loam soils cover approximately 10 percent of the site in the south-central and west-central portions. They are present as the Penn silt loam, three to eight percent slopes, moderately eroded (PeB2 on figure 3.2); the Penn silt loam, three to eight percent slopes, severely eroded (PeB3 on figure 3.2); and the Penn silt loam, eight to 15 percent slopes, moderately eroded (PeC2 on figure 3.2). They are moderately deep to shallow, reddish-brown silt loams that formed in material weathered from red shale, siltstone, and fine-grained sandstone. In a representative profile, the surface layer is friable, reddish-brown silt loam about eight inches thick. The subsoil is reddish-brown shaly silt loam about 12 inches thick, and the substratum is a weak red layer of broken shale pieces with a small amount of silt or loam about 10 to 12 inches thick. Shale, siltstone, or fine-grained sandstone bedrock occurs at a depth of two to three feet. Penn soils have moderately rapid permeability (two to 6.3 inches per hour) and low to moderate available moisture capacity. The soil reaction is strongly acid to neutral (pH, 4.5 to 7.0).<sup>61</sup>

Stony land, steep (StE on figure 3.2) covers approximately 10 percent of the site in a narrow band along the steep slopes of the western portion. This land type occurs in areas of Manor, Penn, Edgemont, and Neshaminy soils; however, individual soil types are not mapped within this unit. These areas are stony and often contain rock ledges and have slopes ranging from 25 to 80 percent. The soil horizons are typically shallow, and the depth to bedrock ranges from several inches to 10 feet. Surface runoff is rapid, and internal drainage is medium to rapid. Permeability and soil reaction values are variable and are not assigned for this unit.<sup>61</sup>

### **3.3.3 Groundwater**

Groundwater in the area occurs under water-table and artesian conditions. Recharge of groundwater is due to the fraction of the local precipitation that infiltrates the soil and underlying material, eventually reaching the saturated zone. From this point, groundwater generally moves downward and laterally until it eventually returns to the surface through discharge points such as springs, seeps, wells, and streambeds.<sup>52,54</sup>

**ORIGINAL**  
**(Red)**

The rocks of the (b) (9) Formation form the aquifer immediately underlying the site. Groundwater in the (b) (9) Formation is mainly stored in and moves through secondary forms of porosity. Due to the fine-grained nature of most of the rocks of the (b) (9) the permeability due to primary porosity is small. Fractures parallel to bedding are usually narrow and probably contribute little to the formation's permeability. The most significant openings with respect to groundwater flow are the nearby vertical joint planes, which cross each other at various angles throughout the beds of the (b) (9) Formation. These provide an interconnected series of channels through which groundwater can flow.<sup>52,54</sup>

The (b) (9) is generally a reliable source of small to moderate supplies of water. Sufficient supplies of water for domestic use can be obtained at almost any location from wells drilled (b) (9) below the water table. In order to obtain maximum yields, wells should be drilled to depths ranging (b) (9) or more. Wells in the (b) (9) Formation of Montgomery County yield from five gpm or less to more than 300 gpm.<sup>52,54</sup>

Forty-two wells in the (b) (9) within Limerick Township range in depth (b) (9) feet, with a median depth (b) (9). These wells have yields ranging from four to 60 gpm, with a median value of 15 gpm. The static water levels in these wells range from (b) (9) the surface, with a median depth of (b) (9) the surface. The median depth to consolidated bedrock is (b) (9).<sup>57</sup>

The (b) (9) is the poorest water-bearing horizon of the (b) (9) of the area; however, domestic supplies are usually available at depths of not more than 250 feet. The rocks have low porosity and low permeability, although joint openings provide secondary porosity and permeability. The least permeable beds are shale and argillite, which have little or no intergranular permeability. Well yields from the (b) (9) in Chester County range from five to 32 gpm, with a median value of 12 gpm.<sup>53,57,59</sup>

The (b) (9) will generally provide sufficient amounts of groundwater for domestic use and will frequently yield adequate supplies for small municipal and industrial uses. Most of the groundwater occurs in and moves through secondary openings along bedding planes, joints, and faults. Domestic wells in the (b) (9) have specific capacities ranging from less than 0.04 to over four gpm per foot of drawdown, with a median value of 0.53 gpm per foot of drawdown. Non-domestic wells have a median-specific capacity of 0.9 gpm per foot of drawdown and a median yield of 120 gpm.<sup>62</sup>

The (b) (9) Formation is capable of yielding moderate to large supplies of water to wells. The more coarse sandstone and conglomerate beds are the best aquifers of the formation. The (b) (9) Formation has high to moderate total effective porosity and permeability resulting from primary intergranular openings in the weathered portion and from secondary joint and bedding-plane openings in the unweathered portion. Well yields in the Stockton Formation of Chester County range from four to 800 gpm, with a median value of 25 gpm.<sup>53,57,59</sup>

The (b) (9) is considered to be a poor formation as a source of water. Most of the groundwater circulation is dependent upon joint openings, which provide very low secondary porosity and low permeability. The (b) (9) is relatively impervious and is expected to act as a barrier to groundwater flow with respect to the adjacent Triassic sedimentary rocks. The (b) (9) will usually yield small amounts of water from relatively shallow depths sufficient for domestic supplies. Well yields in (b) (9) in Chester County range from 0.5 to 40 gpm, with a median value of nine gpm.<sup>53,57,59</sup>

The (b) (9) generally comprises poor water-bearing rocks; however, it yields small supplies sufficient for domestic use. These rocks have low primary porosity, although joints provide secondary porosity and permeability of low magnitude. The best water-bearing zones are in the most weathered and fractured portions above the fresh, consolidated bedrock. Wells in the (b) (9) of Chester County have yields ranging from one to 180 gpm, with a median value of 24 gpm.<sup>53,57,59</sup>

Available records indicate that six industrial supply wells were completed at the site in 1934 and 1935. These wells range in depth (b) (9). All are completed in the (b) (9) Formation and have yields ranging from 10 to 225 gpm, with an average yield of 117.5 gpm. The static water levels and depth of casing in the wells are not reported. These wells were not observed during the site inspection, and their current status is unknown.<sup>4,53</sup>

The actual depth to groundwater and direction of shallow groundwater flow at the site are unknown. Based on topographic control and on the role of rivers and streams as groundwater discharge points, the direction of groundwater flow throughout most of the site is expected to be to the west, southwest, or south toward the Schuylkill River. Based on the site's topographic elevations with respect to the Schuylkill River, the depth to groundwater is expected to be within a few feet of the surface adjacent to the river and to be less than or equal to about 70 feet below the ground surface at the highest elevation of the site. Median water depths in area wells suggest that the depth to groundwater is probably on the order of 40 feet below the ground surface.<sup>1,47,52,54</sup>

### **3.4 CLIMATE AND METEOROLOGY**

According to the climatological data obtained for the city of Phoenixville, Pennsylvania, located approximately six miles southeast of the site, and based on the period from 1951 to 1980, the following is offered. The average annual temperature is 53.1°F; the coldest month is January with a mean temperature of 30.1°F, and the warmest month is July with a mean temperature of 74.9°F. The average annual precipitation is 43.55 inches. The month of highest average precipitation is July, with 4.13 inches; the month with the lowest average precipitation is February, with 2.94 inches. A one-year, 24-hour rainfall will produce 2.5 inches of rain. A two-year, 24-hour rainfall will produce 3.5 inches of rain. The mean annual lake evaporation for the area is 32 inches. The net moisture gain is 11-1/2 inches.<sup>63,64,65,66</sup>

### **3.5 LAND USE**

The site is located in a rural, mostly residential area. State Game Lands No. 234 is to the east. The Schuylkill River borders the property on the southern and western banks. The residential town of Linfield is located adjacent to and north-northeast of the site. ConRail Corporation has abandoned tracks on site.<sup>1,2</sup>

### **3.6 POPULATION DISTRIBUTION**

The Linfield Industrial Park Site is located in the rural outskirts south-southwest of Linfield, Pennsylvania. The population is as follows:<sup>1,34,37</sup>

<u>Distance</u>	<u>Population</u>
0 to 1/4 mile .....	126
1/4 to 1/2 mile .....	338
1/2 to 1 mile .....	1,251
1 to 2 miles .....	9,348
2 to 3 miles .....	6,715
3 to 4 miles .....	15,702
Total .....	34,010

Population figures are based on a count of houses in the study area multiplied by 2.79 persons per house and water supply population figures furnished by CUHWC and PMWA.<sup>1,34,37</sup>

### 3.7 CRITICAL ENVIRONMENTS

Two federally listed endangered birds are expected to be found as transient species in the study area. They are the bald eagle (Haliaeetus leucocephalus) and the peregrine falcon (Falco peregrinus). There is no listed critical habitat for these species in the study area.<sup>67</sup>

## **SECTION 4.0**



#### 4.0 WASTE TYPES AND QUANTITIES

Linfield Industrial Park was operated as a whiskey distillery and a bonded warehouse for the storage and distribution of whiskey products. For an unknown period of time, ethylene glycol (antifreeze) and ammonia-based cleaners were bottled and distributed from the site.<sup>2,13,19</sup> The site has remained inactive since it was sold to 888 Warehousing in 1986. The condition of the site has remained the same since its closure.<sup>2</sup>

There are no records detailing the waste types and quantities generated at the Linfield Industrial Park; however, during the FIT preliminary assessment, numerous items of concern were observed:<sup>2</sup>

##### Parking Lot Area

An abandoned tanker truck was observed northwest of the on-site parking lot. The truck had an approximately 8,000-gallon capacity. The truck's contents were not known.<sup>2</sup>

##### Distillery Area

The distillery building contains approximately eight wooden vats (10 by 20 feet in size) that have an approximately 11,700-gallon capacity. An underground fuel tank with a pump in the distillery area had an approximate capacity of 1,000 gallons. The contents of the fuel tanks are not known. North and west of the distillery building are five above-ground storage tanks. These tanks are metal; the contents are not known. Three of the cylinders are six by 25 feet in size with a 5,284-gallon capacity; two of the cylinders are five by 10 feet in size with a capacity of 939 gallons each. North of these tanks are two underground storage tanks with an approximate capacity of 10,000 gallons each; an unknown quantity of an oil-like substance was inside the tanks. Adjacent to and east of the tanks are an electrical transformer unit and an area of stained soils. The unit measures four by five by four feet in size and contains liquid. Ten 55-gallon drums containing waste oils and unknown materials were observed in this area. Southwest of the distillery building in a shed are ten 55-gallon drums marked "Oil and Varnish." The contents of the drums are not known. South of the distillery building is a metal cylinder above-ground storage tank, 30 by 50 feet in size, with an approximately 264,231-gallon capacity; the contents of the tank are not known.<sup>2</sup>

#### Building No. 2 Area

Located east of building no. 2 are 10 above-ground black metal cylinder storage tanks, six by 40 feet in size, with 8,455-gallon capacities. There are also one white metal cylinder above-ground storage tank, five by 30 feet in size with a 4,400-gallon capacity, and one aluminum-brushed metal cylinder above-ground storage tank, five by 10 feet in size with a 1,467-gallon capacity. Four abandoned tanker trucks were observed around building no. 2. One truck was still hooked to the building via a discharge pipe. The trucks have a capacity of 8,000 gallons each; the contents of the trucks were not known.<sup>2</sup>

#### Warehouse Area

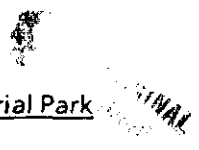
Two aluminum cylinder above-ground storage tanks, 40 by 50 feet in size with a 469,244-gallon capacity, and one aluminum cylinder above-ground storage tank, eight by 15 feet in size with a 5,637-gallon capacity, were located centrally among the warehouse buildings.<sup>2</sup>

#### Concrete Pad Area

At least twenty-two 55-gallon drums were observed in this area: ten 55-gallon drums were empty and rusted, and twelve 55-gallon drums contained unknown liquids. Four drums are located on or near the northwestern pad area, and eight drums are located on or near the southeastern concrete pad area. One 55-gallon drum near the southeastern concrete pad area contained an oil-like residue that had spilled. A stained soil area measuring approximately three by six feet in size was observed in this area. Five 5-gallon metal drums marked "Herbicide" were observed in a shed-like building. A 20-gallon transformer is located south of the pads with stained soils.<sup>2</sup>

#### Former Tank Farm Area

This area once housed five metal cylinder above-ground storage tanks. The tanks were individually fenced and bermed. At an unknown date, the tanks were dismantled and removed. Currently, only the fenced concrete containment areas exist.<sup>1,2,4</sup>



#### Miscellaneous Waste Areas

An unlined limestone pit, adjacent to warehouse no. 10, allegedly was used to dispose of deionized wastewater via groundwater after neutralization. The site, condition, and location of this pit are not known. Also, a catch basin was reportedly utilized for settling suspended solids to reduce the amount of wastewater sent to the wastewater treatment plant. According to PA DER file information, wastewater generated at the site was disposed in the Schuylkill River via the NPDES outfall. Ethylene glycol that spilled into the parking lot drainage system resulted in a contractual removal assistance by Eldridge Waste Management, of Chester, Pennsylvania. The quantity of waste removed is unknown. The soil area at the bottle-making room was contaminated with unknown substances and disposed at the Boyertown Landfill. The quantity of contaminated soils is unknown.<sup>7,8,10,15,17,19,21,22,29</sup>

During an emergency assessment of the site performed by EPA Emergency Response on March 14 and 27, 1991, numerous drums and containers were found within the on-site buildings. These included containers of sulfuric acid, drums and containers marked "Flammable," containers of liquid caustic materials, thousands of small containers filled with antifreeze, drums of methanol, and a drum marked "Toluene."<sup>24,25</sup>

On-site surface soil samples collected by HALLIBURTON NUS ARCS III on January 30, 1992 contained elevated levels of a number of organic and inorganic contaminants, including fluoranthene (72,000 ppb), benzo(a)pyrene (30,000 ppb), PCBs (300,000 ppb), and lead (4,810 ppm).<sup>4</sup>

During the site inspection, a trash dump measuring approximately 30 by 75 feet and consisting mostly of empty plastic antifreeze containers was observed between the southernmost warehouse and the Schuylkill River. Adjacent to this area were approximately 60 upright empty and rusted drums. No markings were visible on the drums.<sup>4</sup>

## **SECTION 5.0**

## 5.0 FIELD TRIP REPORT

### 5.1 SUMMARY

On Tuesday, January 28, 1992, HALLIBURTON NUS ARCS III personnel (b) (6) conducted a site inspection of the Linfield Industrial Park Site in Linfield, Montgomery County, Pennsylvania. Weather conditions were clear with light winds, and temperatures were in the 30s.

Eleven aqueous and 14 solid samples were obtained (see figures 5.1 and 5.2, pages 5-4 and 5-5). Photographs were taken on site (see figure 5.3, page 5-7, and the photograph log, section 5.5).

### 5.2 PERSONS CONTACTED

#### 5.2.1 Prior to Field Trip

Bernard Shafran  
Frenkel and Hershkowitz, P.C.  
Attorneys at Law  
319 Fifth Avenue  
New York, NY 10016  
(212) 679-4666

Zelma Maldonado  
U.S. EPA  
841 Chestnut Building  
Philadelphia, PA 19107  
(215) 597-8333

George Danyliw  
PA DER  
Southeast Regional Office  
Lee Park, Suite 6010  
555 North Lane  
Conshohocken, PA 19428  
(215) 832-6145

Marty Powell  
U.S. EPA  
841 Chestnut Building  
Philadelphia, PA 19107  
(215) 597-6680

#### 5.2.2 At the Site

Bernard Shafran  
Frenkel and Hershkowitz, P.C.  
Attorneys at Law  
319 Fifth Avenue  
New York, NY 10016  
(212) 679-4666

Zelma Maldonado  
U.S. EPA  
841 Chestnut Building  
Philadelphia, PA 19107  
(215) 597-8333

**ORIGINAL  
(Red)**

Site Name: Linfield Industrial Park  
Project No.: 3263-09

**5.2.3 Water Supply Well Information**

The following off-site wells were sampled during the site inspection. (b) (6)

Completed well questionnaires are located in appendix C.

(b) (6)  
(b) (6)  
Linfield, Pennsylvania 19468  
(b) (6)  
Drinking  
(b) (6)

(b) (6)  
(b) (6)  
Linfield, Pennsylvania 19468  
(b) (6)  
Drinking  
(b) (6)

TDD NUMBER 3263-09  
EPA NUMBER PA-2898

### 5.3 SAMPLE LOG

SITE NAME Linfield Industrial Park

TRAFFIC REPORTS			SAMPLE IDENTIFIER	PHASE	SAMPLE DESCRIPTION	SAMPLE LOCATION	TARGET USE	pH	FIELD MEASUREMENTS	
Organic	Inorganic	High Hazard								
CHY 02	MCJE 02		(b) (9)	AQ	Clear, no odor	(b) (9)	Drinking water for the people of Spring City and Poyersford		No HNU readings above bkgd.	
CHY 03	MCJE 03		(b) (9)	AQ	clear, no odor	(b) (9) Linfield, PA	Drinking			
CHY 04	MCJE 04		(b) (9)	AQ	clear, no odor	(b) (9) Linfield, PA	Drinking			
CHY 05	MCJE 05		SW-1	AQ	clear, fish-like odor	Upstream sample from Schuylkill R., under Main St. bridge	Recreation; three water companies operate intakes within 15 miles downstream	6.92	No HNU readings above background	
CHY 06	MCJE 06		SW-2	AQ	clear, no odor	Mid-stream sample from Schuylkill R., at former NPDES discharge	Recreation; three water companies operate water intakes within 15 miles downstream	6.72		
CHY 07	MCJE 07		SW-3	AQ	clear, no odor	On-site drainage swale ~ 30' from confluence with Schuylkill R.	No uses	6.70		
CHY 08	MCJE 08		SW-4	AQ	slightly turbid, no odor	Downstream sample from Schuylkill R.	Recreation; three water companies operate intakes within 15 miles downstream	6.47		
CHY 10	MCJE 10		SW-6	AQ	Duplicate of SW-4	Same as SW-4	same as SW-4			
CHY 09	MCJE 09		(b) (9)	AQ	clear, no odor	(b) (9)	Drinking water for Spring City and Poyersford			

ORIGINAL  
(Red)

TOD NUMBER 3263-09  
EPA NUMBER PA-2898

### 5.3 SAMPLE LOG

**SITE NAME** Linfield Industrial Park

[illegible]



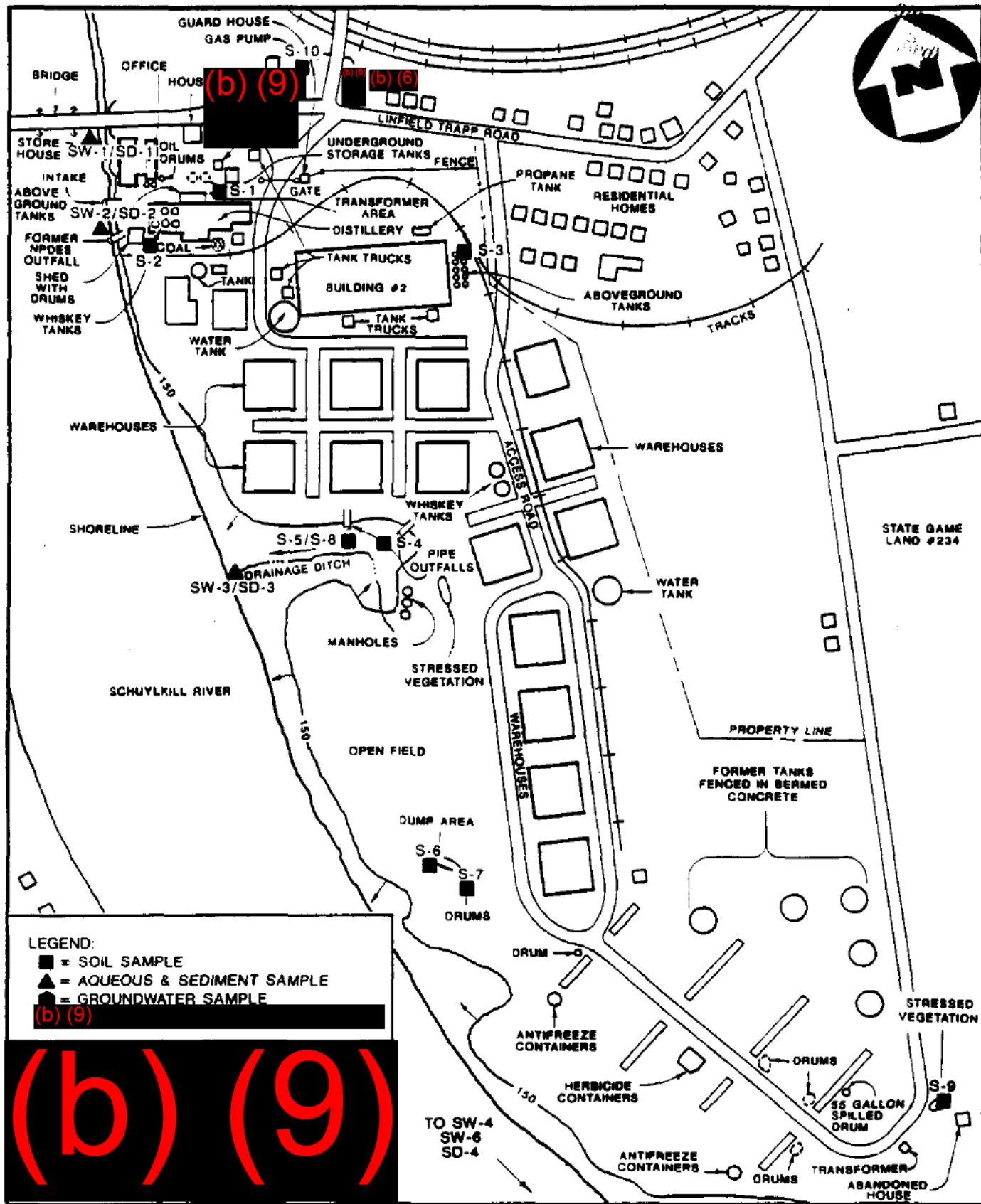
TDD NUMBER 3263-09  
 EPA NUMBER PA-2898

### 5.3 SAMPLE LOG

SITE NAME Linfield Industrial Park

TRAFFIC REPORTS			SAMPLE IDENTIFIER	PHASE	SAMPLE DESCRIPTION	SAMPLE LOCATION	TARGET USE	pH	FIELD MEASUREMENT:	
Organic	Inorganic	High Hazard								
CHY 24	MCJE 24		AQ-0	AQ	Aqueous blank	_____	_____		_____	
CHY 11	MCJE 11		Sd-1	SOL	light brown silt, red shale pebbles, some organic matter	same as SW-1	Access unrestricted		No HNU readings above background	
CHY 12	MCJE 12		Sd-2	SOL	light brown silt, some organic matter	same as SW-2	Access unrestricted			
CHY 13	MCJE 13		Sd-3	SOL	clayey silt, reddish	same as SW-3	Access unrestricted			
CHY 14	MCJE 14		Sd-4	SOL	Brown sandy loam	same as SW-4	Access unrestricted			
CHY 15	MCJE 15		S-1	SOL	dark brown sandy loam on site	At base of transformer on northern side of distillery	Access unrestricted			
CHY 16	MCJE 16		S-2	SOL	black viscous silty clay with grease odor on site	At drums outside of shed west of distillery	Access unrestricted			
CHY 17	MCJE 17		S-3	SOL	Brown loam with some roots and reddish clay on site	Inside above ground tank farm on eastern side of building #2	Access unrestricted			
CHY 18	MCJE 18		S-4	SOL	Brown sandy loam on site	At confluence of two concrete drainage troughs south and west of warehouses	Access unrestricted			

DATE/TIME  
10/1/88  
10:00 AM



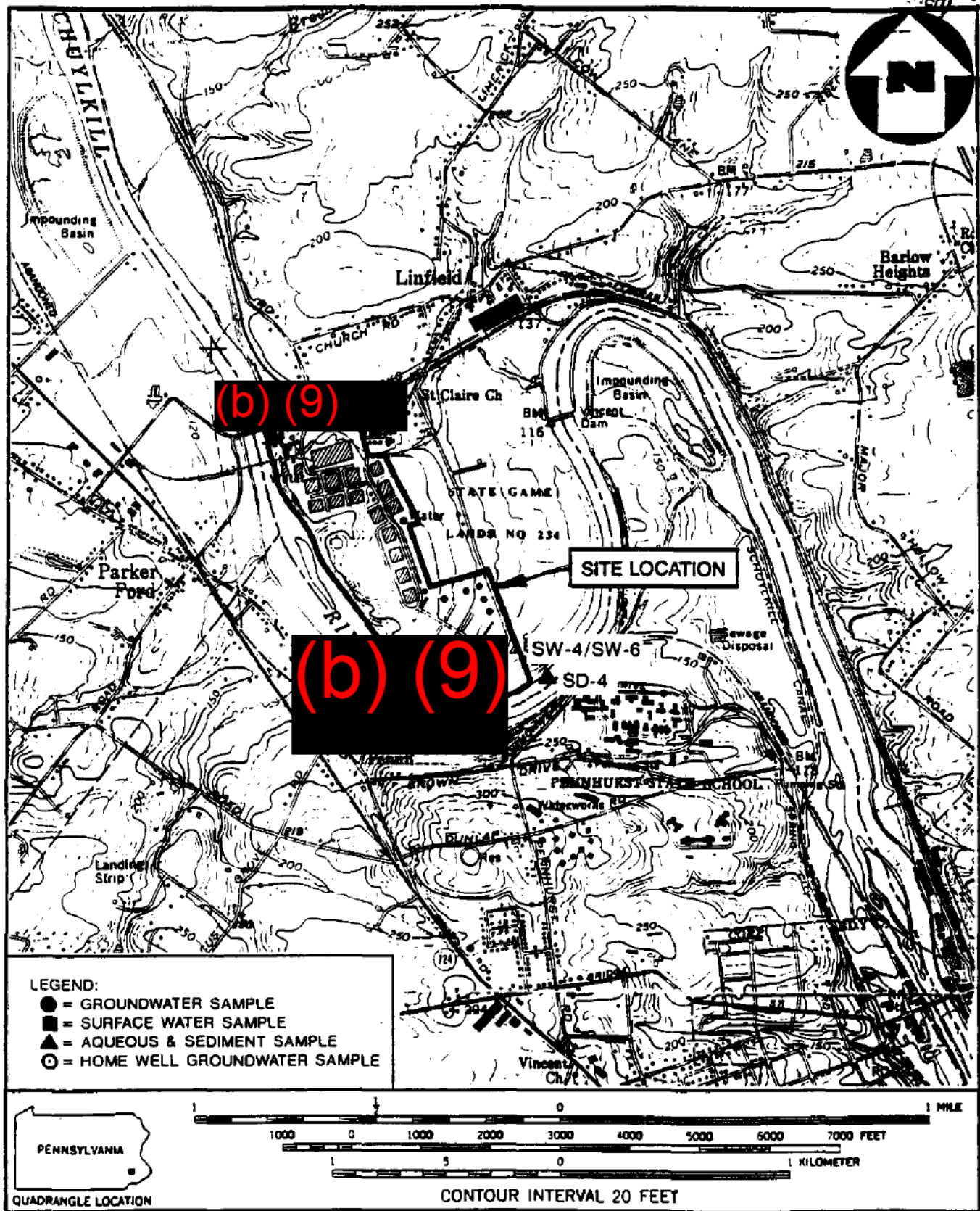
SAMPLE LOCATION MAP

LINFIELD INDUSTRIAL PARK, LINFIELD, PA.

(NO SCALE)

FIGURE 5.1





**OFF-SITE SAMPLE LOCATION MAP**  
**LINFIELD INDUSTRIAL PARK, LINFIELD, PA.**

FIGURE 5.2



#### 5.4 SITE OBSERVATIONS

- The background HNU reading was 0.4 ppm. No readings above background were recorded.
- The radiation mini-alert was set on the X1 position. No readings above background were recorded.
- The entrance gate to the site was fitted with a lock and chain.
- There was a fence on the northeastern corner of the site. Otherwise, pedestrian access to the site was unrestricted.
- The on-site production well was not accessible for sampling. An inoperative pump was inside the well.
- A trash dump, consisting mostly of empty plastic antifreeze containers, was west of the southernmost warehouse. The dump measured approximately 30 by 75 feet. Adjacent to the dump were approximately 60 drums that were upright, empty, and rusted. No markings were visible on the drums.
- Two areas of stressed vegetation were observed on site: an approximately 20- by 100-foot area in the middle of the site, between the warehouse and the Schuylkill River, and an approximately 100- by 100-foot area at the southern end of the site. The vegetation in these two areas was identical.
- Bicycle tracks were observed in the snow between the third and fourth warehouse from the southernmost warehouse.
- Shotgun shells were observed throughout the site.

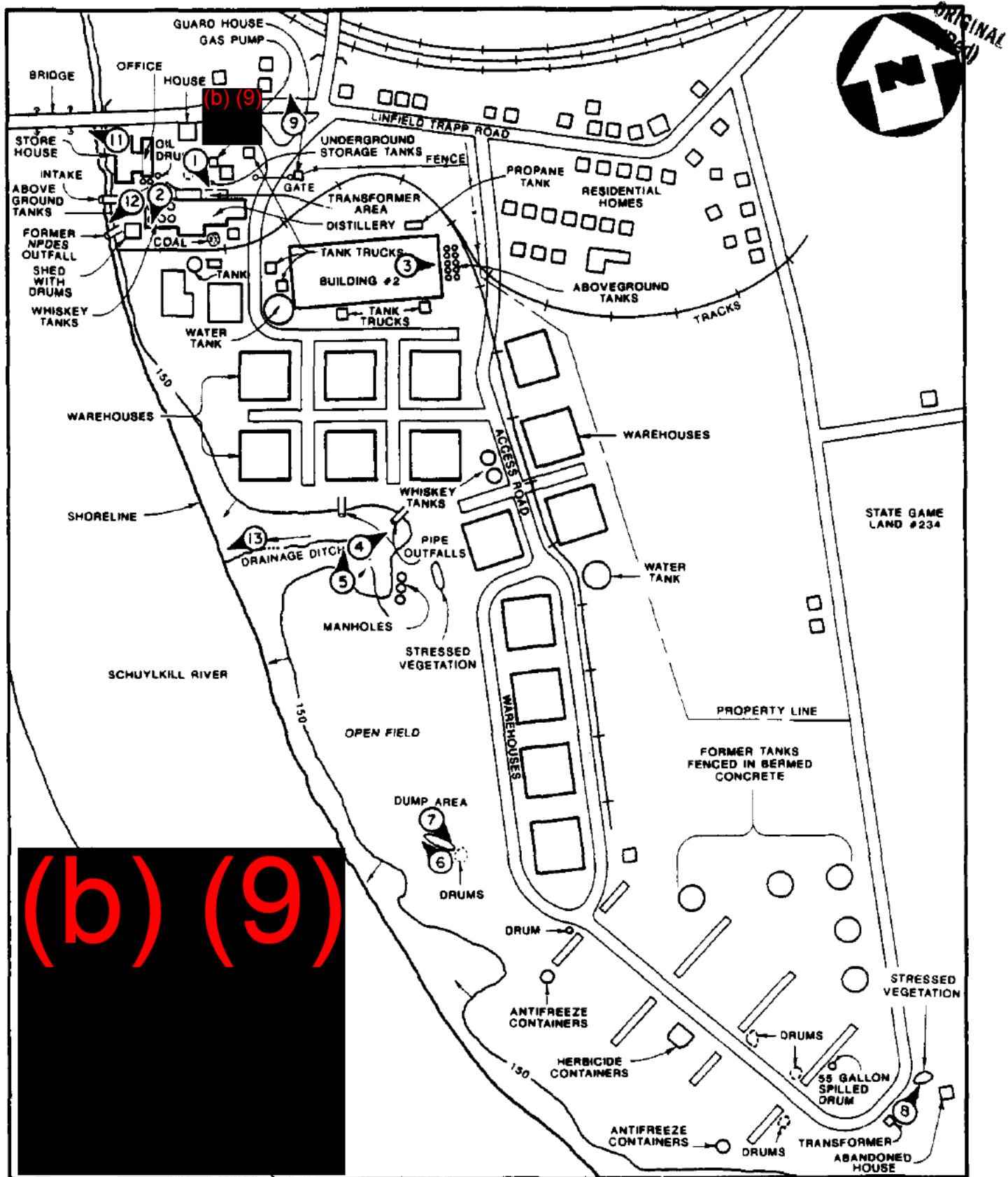


PHOTO LOCATION MAP

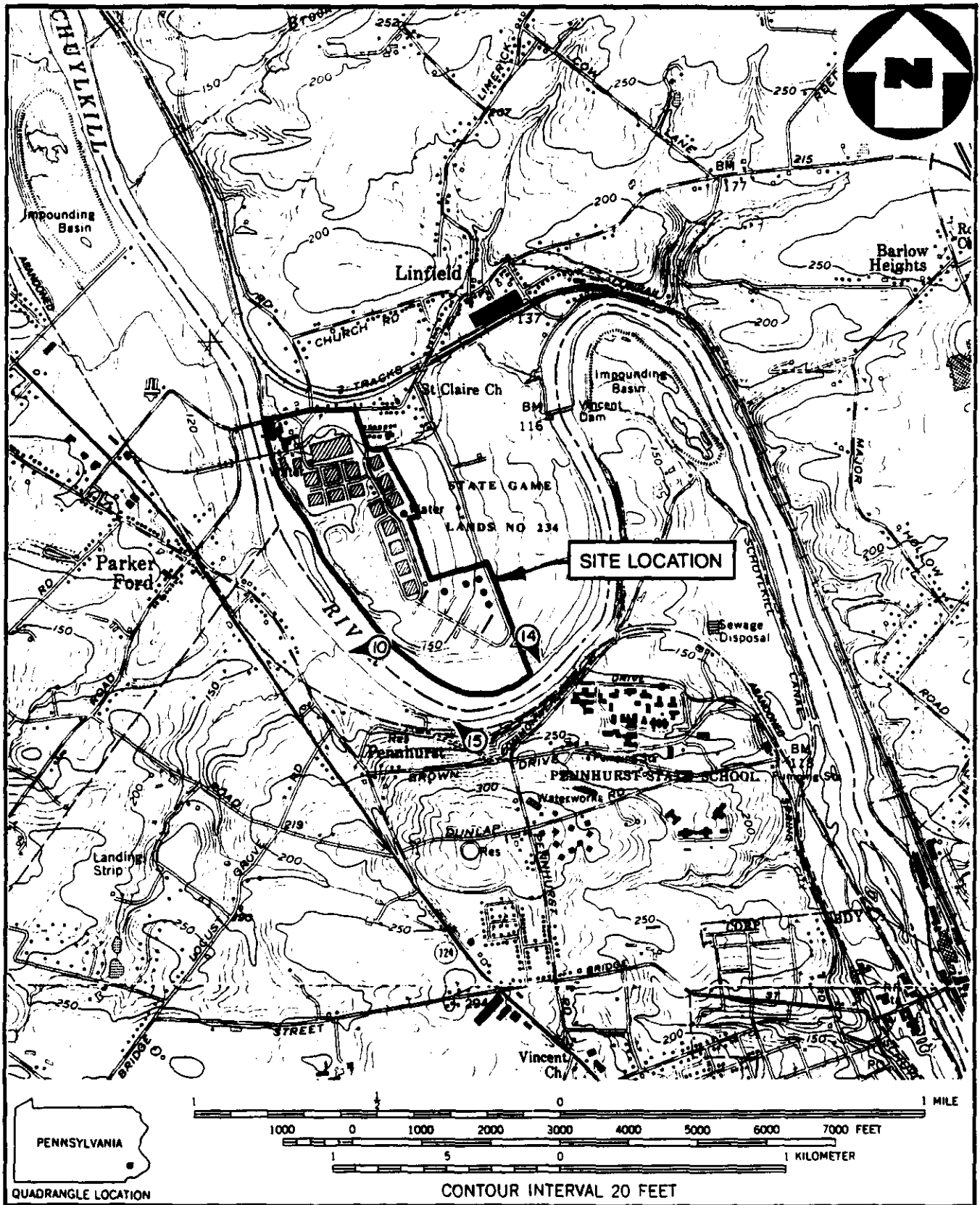
LINFIELD INDUSTRIAL PARK, LINFIELD, PA.

(NO SCALE)

FIGURE 5.3



ORIGINAL  
Red)



SOURCE: (7.5 MINUTE SERIES) U.S.G.S. PHOENIXVILLE, PA QUAD.

**OFF-SITE PHOTO LOCATION MAP**  
**LINFIELD INDUSTRIAL PARK, LINFIELD, PA.**

SCALE 1: 24000

FIGURE 5.4





EPA REGION III  
SUPERFUND DOCUMENT MANAGEMENT SYSTEM

DOC ID # 404788  
PAGE #           

IMAGERY COVER SHEET  
UNSCANNABLE ITEM

Contact the CERCLA Records Center to view this document.

SITE NAME Linfield Ind Park  
OPERABLE UNIT 00  
SECTION/BOX/FOLDER 1C Box 1 1008

REPORT OR DOCUMENT TITLE Final Screening Site  
Inspection  
DATE OF DOCUMENT Sept. 1, 1992  
DESCRIPTION OF IMAGERY photos  
NUMBER AND TYPE OF IMAGERY ITEM(S) 15 photos



EPA

**POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT**  
**PART 1 - SITE LOCATION AND INSPECTION INFORMATION**

**I. IDENTIFICATION**01 STATE  
PA02 SITE NUMBER  
2898**II. SITE NAME AND LOCATION**

01 SITE NAME (Legal, common, or descriptive name of site)

Linfield Industrial Park

02 STREET, ROUTE NO., OR SPECIFIC LOCATION IDENTIFIER

Linfield - Trappe Road

03 CITY

Linfield

04 STATE

PA

05 ZIP CODE

19468

06 COUNTY

Montgomery

07 COUNTY  
CODE

091

08 CONG.  
DIST

PA10

09 COORDINATES

LATITUDE

40° 12' 08" N

LONGITUDE

75° 34' 38" W

10 TYPE OF OWNERSHIP (Check one)

☒ A. PRIVATE☐ B. FEDERAL☐ C. STATE☐ D. COUNTY☐ E. MUNICIPAL☐ F. OTHER☐ G. UNKNOWN**III. INSPECTION INFORMATION**

01 DATE OF INSPECTION

1 / 28 / 92  
month day year

02 SITE STATUS

☐ A. ACTIVE☒ B. INACTIVE

03 YEARS OF OPERATION

unknown / 1986

BEGINNING YEAR

ENDING YEAR

UNKNOWN

04 AGENCY PERFORMING INSPECTION (Check all that apply)

☐ A. EPA☒ B. EPA CONTRACTOR

HALLIBURTON NUS

(Name of firm)

☐ C. MUNICIPAL☐ D. MUNICIPAL CONTRACTOR

(Name of firm)

☐ E. STATE☐ F. STATE CONTRACTOR

(Name of firm)

☐ G. OTHER

(Specify)

05 CHIEF INSPECTOR

(b) (4)

06 TITLE

Geologist

07 ORGANIZATION

HALLIBURTON NUS

08 TELEPHONE NO.

(215) 971-0900

09 OTHER INSPECTORS

(b) (4)

10 TITLE

Environmental Scientist

11 ORGANIZATION

HALLIBURTON NUS

12 TELEPHONE NO.

(215) 971-0900

(b) (4)

Environmental Scientist

HALLIBURTON NUS

(215) 971-0900

(b) (4)

Environmental Scientist

HALLIBURTON NUS

(215) 971-0900

(b) (4)

Environmental Scientist

HALLIBURTON NUS

(215) 971-0900

(b) (4)

Environmental Scientist

HALLIBURTON NUS

(215) 971-0900

13 SITE REPRESENTATIVES INTERVIEWED

Bernard Shafran

14 TITLE

Attorney

15 ADDRESS

319 Fifth Avenue  
New York, New York 10016

16 TELEPHONE NO.

(212) 679-4666

17 ACCESS GAINED BY  
(Check one)☒ PERMISSION☐ WARRANT

18 TIME OF INSPECTION

8:30 a.m.

19 WEATHER CONDITIONS

Clear and cold

**IV. INFORMATION AVAILABLE FROM**

01 CONTACT

Michael Giuranna

02 OF (Agency/Organization)

United States  
Environmental Protection Agency

03 TELEPHONE NO.

(215) 597-3165

04 PERSON RESPONSIBLE FOR SITE INSPECTION FORM

(b) (4)

05 AGENCY

HALLIBURTON  
NUS

06 ORGANIZATION

ARCS III

07 TELEPHONE NO.

(215) 971-0900

08 DATE

2 / 12 / 92  
month day year





ORIGINAL  
EPA

POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 2 - WASTE INFORMATION

I. IDENTIFICATION

01 STATE  
PA

02 SITE NUMBER  
2898

II. WASTE STATES, QUANTITIES, AND CHARACTERISTICS

01 PHYSICAL STATES (Check all that apply)

- ☐ A. SOLID  
☐ B. POWDER, FINES  
☐ C. SLUDGE  
☐ D. OTHER \_\_\_\_\_  
(Specify)  
☐ E. SLURRY  
☒ F. LIQUID  
☐ G. GAS

02 WASTE QUANTITY AT SITE

(Measures of waste quantities must be independent)

TONS unknown  
CUBIC YARDS \_\_\_\_\_  
NO. OF DRUMS \_\_\_\_\_

03 WASTE CHARACTERISTICS (Check all that apply)

- ☒ A. TOXIC  
☒ B. CORROSIVE  
☐ C. RADIOACTIVE  
☒ D. PERSISTENT  
☒ E. SOLUBLE  
☐ F. INFECTIOUS  
☒ G. FLAMMABLE  
☐ H. IGNITABLE  
☐ I. HIGHLY VOLATILE  
☐ J. EXPLOSIVE  
☐ K. REACTIVE  
☐ L. INCOMPATIBLE  
☐ M. NOT APPLICABLE

III. WASTE TYPE

CATEGORY	SUBSTANCE NAME	01 GROSS AMOUNT	02 UNIT OF MEASURE	03 COMMENTS
SLU	SLUDGE			
OLW	OILY WASTES	unknown		On-site soil samples taken by HALLIBURTON NUS ARCS III
SOL	SOLVENTS	unknown		on January 30, 1992 revealed elevated levels of
PSD	PESTICIDES			organic and inorganic contaminants. On March 27, 1991,
OCC	OTHER ORGANIC CHEMICALS	unknown		EPA Emergency Response found containers of various
IOC	INORGANIC CHEMICALS	unknown		hazardous materials, including acids, caustics, and
ACD	ACIDS	unknown		flammables, inside the on-site warehouses.
BAS	BASES	unknown		
MES	HEAVY METALS	unknown		

IV. HAZARDOUS SUBSTANCES (See Appendix for most frequently cited CAS Numbers)

01 CATEGORY	02 SUBSTANCE NAME	03 CAS NUMBER	04 STORAGE DISPOSAL METHOD	05 CONCENTRATION	06 MEASURE OF CONCENTRATION
SOL	toluene	108-88-3	on-site soil	120	ppb
OCC	fluoranthene	206-44-0	on-site soil	72,000	ppb
OCC	pyrene	129-00-0	on-site soil	54,000	ppb
OCC	chrysene	218-01-9	on-site soil	32,000	ppb
OCC	benzo(a)pyrene	50-32-8	on-site soil	30,000	ppb
OCC	phenanthrene	85-01-8	on-site soil	52,000	ppb
OCC	anthracene	120-12-7	on-site soil	14,000	ppb
IOC	Aroclor 1260	11096-82-5	on-site soil	300,000	ppb
MES	lead	7439-92-1	on-site soil	4,810	ppm
MES	mercury	7439-97-6	on-site soil	2.90	ppm
MES	zinc	7440-66-6	on-site soil	1,230	ppm
IOC	cyanide		on-site soil	1.8	ppm

IV. FEEDSTOCKS (See Appendix for CAS Numbers) N/A

CATEGORY	01 FEEDSTOCK NAME	02 CAS NUMBER	CATEGORY	01 FEEDSTOCK NAME	02 CAS NUMBER
FDS			FDS		
FDS			FDS		
FDS			FDS		
FDS			FDS		

VI. SOURCES OF INFORMATION (Cite specific references, e.g., state files, sample analysis, reports)

HALLIBURTON NUS, ARCS III. Site inspection; site visit. Project No. 3263-09, January 30, 1992.  
NUS Corporation, FIT 3. Preliminary assessment. TDD No. F3-9002-18, June 27, 1991.

**EPA**

**POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 3 - DESCRIPTION OF HAZARDOUS CONDITIONS AND INCIDENTS**

**I. IDENTIFICATION**

01 STATE PA	02 SITE NUMBER 2898
----------------	------------------------

**II. HAZARDOUS CONDITIONS AND INCIDENTS**

01 ☒ A. GROUNDWATER CONTAMINATION 02 ☐ OBSERVED (DATE: \_\_\_\_\_) ☒ POTENTIAL ☐ ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: 19,500 04 NARRATIVE DESCRIPTION

On-site soil samples taken by ARCS III on January 30, 1992 contained elevated levels of organic and inorganic contaminants, including benzo(a)pyrene (30 ppm), PCBs (300 ppm), and lead (4,810 ppm). 19,831 people use groundwater within the study area for drinking.

01 ☒ B. SURFACE WATER CONTAMINATION 02 ☒ OBSERVED (DATE: \_\_\_\_\_) ☐ POTENTIAL ☐ ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: 38,894 04 NARRATIVE DESCRIPTION

Aqueous samples from the Schuylkill River taken by ARCS III on January 30, 1992 contained cyanide (11.70 ppb), zinc (301 J ppb), and lead (59.9 J ppb). 38,894 people obtain drinking water from the Schuylkill River within 15 miles downstream from the site.

01 ☐ C. CONTAMINATION OF AIR 02 ☐ OBSERVED (DATE: \_\_\_\_\_) ☐ POTENTIAL ☐ ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: \_\_\_\_\_ 04 NARRATIVE DESCRIPTION

None reported or observed.

01 ☐ D. FIRE/EXPLOSIVE CONDITIONS 02 ☐ OBSERVED (DATE: \_\_\_\_\_) ☐ POTENTIAL ☐ ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: \_\_\_\_\_ 04 NARRATIVE DESCRIPTION

None reported or observed.

01 ☒ E. DIRECT CONTACT 02 ☐ OBSERVED (DATE: \_\_\_\_\_) ☒ POTENTIAL ☐ ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: 1,715/1 mile 04 NARRATIVE DESCRIPTION

On-site soil samples taken by ARCS III on January 30, 1992 contained elevated levels of organic and inorganic contaminants including benzo(a)pyrene (30 ppm), PCBs (300 ppm), and lead (4,810 ppm). Site access is unrestricted. 1,819 people live within one mile of the site.

01 ☒ F. CONTAMINATION OF SOIL 02 ☒ OBSERVED (DATE: \_\_\_\_\_) ☐ POTENTIAL ☐ ALLEGED  
03 AREA POTENTIALLY AFFECTED: 125 (Acres) 04 NARRATIVE DESCRIPTION

On-site soil samples taken by ARCS III on January 30, 1992 contained elevated levels of organic and inorganic contaminants, including benzo(a)pyrene (30 ppm), PCBs (300 ppm), and lead (4,810 ppm). The total area of the site is 125 acres.

01 ☒ G. DRINKING WATER CONTAMINATION 02 ☐ OBSERVED (DATE: \_\_\_\_\_) ☒ POTENTIAL ☐ ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: 58,394 04 NARRATIVE DESCRIPTION

On-site soil samples taken by ARCS III on January 30, 1992 contained elevated levels of organic and inorganic contaminants, and an aqueous sample from the Schuylkill River contained cyanide (11.70 ppb), zinc (301 J ppb), and lead (59.9 J ppb). 58,275 people obtain drinking water from either groundwater within the study area or surface water within 15 miles downstream from the site.

01 ☐ H. WORKER EXPOSURE/INJURY 02 ☐ OBSERVED (DATE: \_\_\_\_\_) ☐ POTENTIAL ☐ ALLEGED  
03 WORKERS POTENTIALLY AFFECTED: \_\_\_\_\_ 04 NARRATIVE DESCRIPTION

None reported or observed.

01 ☒ I. POPULATION EXPOSURE/INJURY 02 ☐ OBSERVED (DATE: \_\_\_\_\_) ☒ POTENTIAL ☐ ALLEGED  
03 POPULATION POTENTIALLY AFFECTED: 58,394 04 NARRATIVE DESCRIPTION

The population potentially affected by the site consists of those obtaining drinking water from companies drawing from the Schuylkill River, those utilizing groundwater drawn from the study area, and those persons accessing the contaminated areas of the site.



ORIGINAL  
EPA (Red)



POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT

PART 3 - DESCRIPTION OF HAZARDOUS CONDITIONS AND INCIDENTS

I. IDENTIFICATION

01 STATE  
PA

02 SITE NUMBER  
2898

II. HAZARDOUS CONDITIONS AND INCIDENTS (Continued)

01 ☐ J. DAMAGE TO FLORA

02 ☐ OBSERVED (DATE: \_\_\_\_\_)

☐ POTENTIAL

☐ ALLEGED

04 NARRATIVE DESCRIPTION

None reported or observed.

01 ☒ K. DAMAGE TO FAUNA

02 ☐ OBSERVED (DATE: \_\_\_\_\_)

☒ POTENTIAL

☐ ALLEGED

04 NARRATIVE DESCRIPTION (Include name(s) of species)

Aqueous samples from the Schuylkill River taken by ARCS III on January 30, 1992 contained cyanide (11.7 ppb), zinc (301 J ppb), and lead (59.9 J ppb). The types of species potentially affected are not known.

01 ☒ L. CONTAMINATION OF FOOD CHAIN

02 ☐ OBSERVED (DATE: \_\_\_\_\_)

☒ POTENTIAL

☐ ALLEGED

04 NARRATIVE DESCRIPTION

Aqueous samples from the Schuylkill River taken by ARCS III on January 30, 1992 contained cyanide (11.7 ppb), zinc (301 J ppb), and lead (59.9 J ppb).

01 ☒ M. UNSTABLE CONTAINMENT OF WASTES

(Spills, Runoff, Standing liquids, Leaking drums)

02 ☒ OBSERVED (DATE: 1/30/92)

☐ POTENTIAL

☐ ALLEGED

03 POPULATION POTENTIALLY AFFECTED: 1,715

04 NARRATIVE DESCRIPTION

On-site soil samples taken by ARCS III on January 30, 1992 contained elevated levels of organic and inorganic contaminants including benzo(a)pyrene (30 ppm), PCBs (300 ppm), and lead (4,810 ppm). Approximately 1,819 people reside within a one-mile radius.

01 ☐ N. DAMAGE TO OFFSITE PROPERTY

02 ☐ OBSERVED (DATE: \_\_\_\_\_)

☐ POTENTIAL

☐ ALLEGED

04 NARRATIVE DESCRIPTION

None reported or observed.

01 ☐ O. CONTAMINATION OF SEWERS, STORM DRAINS, WWTPs

02 ☐ OBSERVED (DATE: \_\_\_\_\_)

☐ POTENTIAL

☐ ALLEGED

04 NARRATIVE DESCRIPTION

None reported or observed.

01 ☒ P. ILLEGAL/UNAUTHORIZED DUMPING

02 ☒ OBSERVED (DATE: 1/30/92)

☐ POTENTIAL

☐ ALLEGED

04 NARRATIVE DESCRIPTION

An on-site trash dump was observed by ARCS III on January 30, 1992.

05 DESCRIPTION OF ANY OTHER KNOWN, POTENTIAL, OR ALLEGED HAZARDS

None reported or observed.

III. TOTAL POPULATION POTENTIALLY AFFECTED: 58,394

IV. COMMENTS

None

V. SOURCES OF INFORMATION (Cite specific references, e.g., state files, sample analysis, reports)

HALLIBURTON NUS ARCS III. Site inspection; site visit. Project No. 3263-09, January 30, 1992.

**EPA****POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 4 - PERMIT AND DESCRIPTIVE INFORMATION****I. IDENTIFICATION**01 STATE  
PA02 SITE NUMBER  
2898**II. PERMIT INFORMATION**

01 TYPE OF PERMIT ISSUED (Check all that apply)	02 PERMIT NUMBER	03 DATE ISSUED	04 EXPIRATION DATE	05 COMMENTS
<input checked="" type="checkbox"/> A. NPDES	PA0013293	unknown	12/23/88	
<input type="checkbox"/> B. UIC				
<input type="checkbox"/> C. AIR				
<input type="checkbox"/> D. RCRA				
<input type="checkbox"/> E. RCRA INTERIM STATUS				
<input type="checkbox"/> F. SPCC PLAN				
<input type="checkbox"/> G. STATE (specify)				
<input type="checkbox"/> H. LOCAL (specify)				
<input type="checkbox"/> I. OTHER (specify)				
<input type="checkbox"/> J. NONE				

**III. SITE DESCRIPTION**

01 STORAGE/DISPOSAL (Check all that apply)	02 AMOUNT	03 UNIT OF MEASURE	04 TREATMENT (Check all that apply)	05 OTHER
<input type="checkbox"/> A. SURFACE IMPOUNDMENT			<input type="checkbox"/> A. INCINERATION	<input checked="" type="checkbox"/> A. BUILDINGS ON SITE
<input type="checkbox"/> B. PILES			<input type="checkbox"/> B. UNDERGROUND INJECTION	
<input checked="" type="checkbox"/> C. DRUMS, ABOVE GROUND	75 to 100	drums	<input type="checkbox"/> C. CHEMICAL/PHYSICAL	
<input checked="" type="checkbox"/> D. TANK, ABOVE GROUND	21	tanks	<input type="checkbox"/> D. BIOLOGICAL	
<input checked="" type="checkbox"/> E. TANK, BELOW GROUND	2	tanks	<input type="checkbox"/> E. WASTE OIL PROCESSING	06 AREA OF SITE
<input type="checkbox"/> F. LANDFILL			<input type="checkbox"/> F. SOLVENT RECOVERY	
<input type="checkbox"/> G. LANDFARM			<input type="checkbox"/> G. OTHER RECYCLING/RECOVERY	
<input checked="" type="checkbox"/> H. OPEN DUMP	unknown		<input type="checkbox"/> H. OTHER (Specify)	125 (Acres)
<input checked="" type="checkbox"/> I. OTHER (Specify) <u>tanker trucks</u>	4			

**07 COMMENTS**

The site, abandoned since 1986, had been used as a whiskey distillery, a whiskey warehousing operation, and an antifreeze and ammonia-based cleaner repackaging operation. NUS FIT 3 recommended action by EPA Emergency Response after observing numerous drums and tanks on site during a preliminary assessment in February 1991.

**IV. CONTAINMENT****01 CONTAINMENT OF WASTES (Check one)**☐ A. ADEQUATE, SECURE ☐ B. MODERATE ☒ C. INADEQUATE, POOR ☐ D. INSECURE, UNSOUND, DANGEROUS**02 DESCRIPTION OF DRUMS, DIKING, LINERS, BARRIERS, ETC.**

No liners exist for on-site soil contamination detected by ARCS III on January 30, 1992.

**V. ACCESSIBILITY**01 WASTE EASILY ACCESSIBLE: ☒ YES ☐ NO**02 COMMENTS**

On-site surface soil contamination was detected by ARCS III on January 30, 1992. Site access is unrestricted.

**VI. SOURCES OF INFORMATION (Cite specific references, e.g., state files, sample analysis, reports)**

NUS Corporation, FIT 3. Preliminary assessment. TOD No. F3-9008-01, June 27, 1991.

EPA file information.

HALLIBURTON NUS ARCS III. Site inspection; site visit. Project No. 3263-09, January 30, 1992.

**EPA ORIGINAL**

(F-001)

**POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 5. WATER, DEMOGRAPHIC, AND ENVIRONMENTAL DATA****I. IDENTIFICATION**01 STATE  
PA02 SITE NUMBER  
2898**II. DRINKING WATER SUPPLY**01 TYPE OF DRINKING SUPPLY  
(Check as applicable)

SURFACE

WELL

ENDANGERED

AFFECTED

MONITORED

COMMUNITY

A. ☒B. ☒A. ☐B. ☐C. ☐

NON-COMMUNITY

C. ☐D. ☒D. ☐E. ☐F. ☐

03 DISTANCE TO SITE

A. Approximately 0.1 (mi)

B. Approximately 100 to 200 feet (mi)

**III. GROUNDWATER**

01 GROUNDWATER USE IN VICINITY (Check one)

☒ A. ONLY SOURCE FOR DRINKING☐ B. DRINKING  
(Other sources available)  
COMMERCIAL, INDUSTRIAL, IRRIGATION  
(No other water sources available)☐ C. COMMERCIAL, INDUSTRIAL, IRRIGATION  
(Limited other sources available)☐ D. NOT USED, UNUSABLE

02 POPULATION SERVED BY GROUND WATER

19,500

03 DISTANCE TO NEAREST DRINKING WATER WELL

100 to 200 feet (mi)

04 DEPTH TO GROUNDWATER

Estimated  
0 to 70 (ft)

05 DIRECTION OF GROUNDWATER FLOW

west through south

06 DEPTH TO AQUIFER  
OF CONCERNEstimated  
0 to 70 (ft)07 POTENTIAL YIELD  
OF AQUIFER

432,000 (gpd)

08 SOLE SOURCE AQUIFER

☐ YES ☒ NO

09 DESCRIPTION OF WELLS (Including usage, depth, and location relative to population and buildings)

(b) (9)

10 RECHARGE AREA

☒ YES☐ NOCOMMENTS: Groundwater recharge from  
precipitation that infiltrates site soils  
and rock to saturated zone.

11 DISCHARGE AREA

☒ YES☐ NO

COMMENTS

Expected groundwater discharge to Schuylkill River

**IV. SURFACE WATER**

01 SURFACE WATER USE IN VICINITY (Check one)

☒ A. RESERVOIR, RECREATION  
DRINKING WATER SOURCE☐ B. IRRIGATION, ECONOMICALLY  
IMPORTANT RESOURCES☐ C. COMMERCIAL, INDUSTRIAL☐ D. NOT CURRENTLY USED

02 AFFECTED/POTENTIALLY AFFECTED BODIES OF WATER

NAME:

Schuylkill River

AFFECTED

☒☐☐

DISTANCE TO SITE

0 (mi)

(mi)

(mi)

**V. DEMOGRAPHIC AND PROPERTY INFORMATION**

01 TOTAL POPULATION WITHIN

ONE (1) MILE OF SITE

TWO (2) MILES OF SITE

THREE (3) MILES OF SITE

A. 1,819  
NO. OF PERSONSB. 12,965  
NO. OF PERSONSC. 18,308  
NO. OF PERSONS

02 DISTANCE TO NEAREST POPULATION

&lt; 0.1 (mi)

03 NUMBER OF BUILDINGS WITHIN TWO (2) MILES OF SITE

4,647

04 DISTANCE TO NEAREST OFF-SITE BUILDING

&lt; 0.1 (mi)

05 POPULATION WITHIN VICINITY OF SITE (Provide narrative description of nature of population within vicinity of site, e.g., rural, village, densely populated urban area)

The residential town of Linfield is adjacent to and north-northeast of the site. The residential towns of Parker Ford and Pennhurst are across the Schuylkill River from the site to the west and south, respectively.

**EPA**

**POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 5 - WATER, DEMOGRAPHIC, AND ENVIRONMENTAL DATA**

**I. IDENTIFICATION**01 STATE  
PA02 SITE NUMBER  
2898**VI. ENVIRONMENTAL INFORMATION**

## -01 PERMEABILITY OF UNSATURATED ZONE (Check one)

☐ A.  $10^{-6} - 10^{-8}$  cm/sec    ☐ B.  $10^{-4} - 10^{-6}$  cm/sec    ☐ C.  $10^{-4} - 10^{-3}$  cm/sec    ☒ D. GREATER THAN  $10^{-3}$  cm/sec

## 02 PERMEABILITY OF BEDROCK (Check one)

☐ A. IMPERMEABLE  
(Less than  $10^{-6}$  cm/sec)    ☐ B. RELATIVELY IMPERMEABLE  
( $10^{-6} - 10^{-5}$  cm/sec)    ☒ C. RELATIVELY PERMEABLE  
( $10^{-5} - 10^{-4}$  cm/sec)    ☐ D. VERY PERMEABLE  
(greater than  $10^{-4}$  cm/sec)03 DEPTH TO BEDROCK  
Approximately  
0 to 15 (ft)04 DEPTH OF CONTAMINATED SOIL ZONE  
unknown (ft)05 SOIL pH  
4.5 to 7.0

## 06 NET PRECIPITATION

11.5 (in)

## 07 ONE-YEAR 24-HOUR RAINFALL

2.5 (in)

## 08 SLOPE

SITE SLOPE

10 %

DIRECTION OF SITE SLOPE

southwest

TERRAIN AVERAGE SLOPE

10 %

## 09 FLOOD POTENTIAL

SITE IS IN 500 YEAR FLOOD PLAIN

10 N/A

☐ SITE IS ON BARRIER ISLAND, COASTAL HIGH HAZARD AREA, RIVERINE FLOODWAY

## 11 DISTANCE TO WETLANDS (5-acre minimum)

ESTUARINE

OTHER

A. N/A (mi)

B. 0.1 (mi)

## 12 DISTANCE TO CRITICAL HABITAT (of endangered species)

N/A (mi)

ENDANGERED SPECIES:

## 13 LAND USE IN VICINITY

DISTANCE TO:

COMMERCIAL/INDUSTRIAL

A. 0.5 (mi)

RESIDENTIAL AREAS: NATIONAL/STATE PARKS,  
FORESTS, OR WILDLIFE RESERVES

B. &lt; 0.1 (mi)

PRIME AG LAND

C. unknown (mi)

AGRICULTURAL LANDS

AG LAND

D. (mi)

## 14 DESCRIPTION OF SITE IN RELATION TO SURROUNDING TOPOGRAPHY

The site lies within a meander of the Schuylkill River, which borders the site to the west and south and is approximately 2,500 feet to the east.

**VII. SOURCES OF INFORMATION** (Cite specific references, e.g., state files, sample analysis, reports)

NUS Corporation, FIT 3. Preliminary assessment. TDD No. F3-9008-01, June 27, 1991.

**EPA**

ORIGINAL

(Red)

**POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 6 - SAMPLE AND FIELD INFORMATION****I. IDENTIFICATION**01 STATE  
PA02 SITE NUMBER  
2898**II. SAMPLES TAKEN**

SAMPLE TYPE	01 NUMBER OF SAMPLES TAKEN	02 SAMPLES SENT TO	03 ESTIMATED DATE RESULTS AVAILABLE
GROUNDWATER	3	Organic: Compuchem Inorganic: Keytx	current
SURFACE WATER	6	Organic: Compuchem Inorganic: Keytx	current
WASTE			
AIR			
RUNOFF			
SPILL			
SOIL	14	Organic: Compuchem Inorganic: Keytx	current
VEGETATION			
OTHER			

**III. FIELD MEASUREMENTS TAKEN**

01 TYPE	02 COMMENTS
pH	Surface water pH levels ranged from 6.70 to 6.97.
HNU	No readings above background were recorded.
Mini-Alert	No readings above background were recorded.

**IV. PHOTOGRAPHS AND MAPS**

01 TYPE	<input checked="" type="checkbox"/> GROUND <input type="checkbox"/> AERIAL	02 IN CUSTODY OF <u>U.S. EPA</u> <small>(Name of organization or individual)</small>
03 MAPS <input checked="" type="checkbox"/> YES <input type="checkbox"/> NO	04 LOCATION OF MAPS <u>HALLIBURTON NUS ARCS III</u>	

**V. OTHER FIELD DATA COLLECTED** (Provide narrative description)

N/A

**VI. SOURCES OF INFORMATION** (Cite specific references, e.g., state files, sample analysis, reports)

HALLIBURTON NUS ARCS III. Site inspection; site visit. Project No. 3263-09, January 30, 1992.

**EPA****POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 7 - OWNER INFORMATION****I. IDENTIFICATION**01 STATE  
PA02 SITE NUMBER  
2898**II. CURRENT OWNER(S)****PARENT COMPANY (if applicable)**

01 NAME 888 Warehousing, Incorporated			02 D & B NUMBER			10 NAME			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.) 527 Redford Avenue			04 SIC CODE			12 STREET ADDRESS (P.O. Box, RFD #, Etc.)			13 SIC CODE		
05 CITY Brooklyn		06 STATE NY	07 ZIP CODE 11211			14 CITY		15 STATE	16 ZIP CODE		
01 NAME			02 D & B NUMBER			10 NAME			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)			04 SIC CODE			12 STREET ADDRESS (P.O. Box, RFD #, Etc.)			13 SIC CODE		
05 CITY		06 STATE	07 ZIP CODE			14 CITY		15 STATE	16 ZIP CODE		
01 NAME			02 D & B NUMBER			10 NAME			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)			04 SIC CODE			12 STREET ADDRESS (P.O. Box, RFD #, Etc.)			13 SIC CODE		
05 CITY		06 STATE	07 ZIP CODE			14 CITY		15 STATE	16 ZIP CODE		
01 NAME			02 D & B NUMBER			10 NAME			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)			04 SIC CODE			12 STREET ADDRESS (P.O. Box, RFD #, Etc.)			13 SIC CODE		
05 CITY		06 STATE	07 ZIP CODE			14 CITY		15 STATE	16 ZIP CODE		
01 NAME			02 D & B NUMBER			10 NAME			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)			04 SIC CODE			12 STREET ADDRESS (P.O. Box, RFD #, Etc.)			13 SIC CODE		
05 CITY		06 STATE	07 ZIP CODE			14 CITY		15 STATE	16 ZIP CODE		

**III. PREVIOUS OWNERS(S) (List most recent first)****IV. REALTY OWNER(S) (if applicable, list most recent first)**

01 NAME Publicker Industries, Incorporated			02 D & B NUMBER			10 NAME			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.) 777 West Putnam Avenue			04 SIC CODE			12 STREET ADDRESS (P.O. Box, RFD #, Etc.)			13 SIC CODE		
05 CITY Greenwich		06 STATE CT	07 ZIP CODE 06836			14 CITY		15 STATE	16 ZIP CODE		
01 NAME			02 D & B NUMBER			10 NAME			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)			04 SIC CODE			12 STREET ADDRESS (P.O. Box, RFD #, Etc.)			13 SIC CODE		
05 CITY		06 STATE	07 ZIP CODE			14 CITY		15 STATE	16 ZIP CODE		
01 NAME			02 D & B NUMBER			10 NAME			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)			04 SIC CODE			12 STREET ADDRESS (P.O. Box, RFD #, Etc.)			13 SIC CODE		
05 CITY		06 STATE	07 ZIP CODE			14 CITY		15 STATE	16 ZIP CODE		

**IV. SOURCES OF INFORMATION (Cite specific references, e.g., state files, sample analysis, reports)**

NUS Corporation, FIT 3. Preliminary assessment. TDD No. F3-9008-01, June 27, 1991.



**EPA****ORIGINAL  
(Red)****POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 8 - OPERATOR INFORMATION****I. IDENTIFICATION**01 STATE  
PA02 SITE NUMBER  
2898**II. CURRENT OPERATOR** *(Provide if different from owner)***OPERATOR'S PARENT COMPANY** *(if applicable)*

01 NAME N/A			02 D & B NUMBER			10 NAME N/A			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)				04 SIC CODE		12 STREET ADDRESS (P.O. Box, RFD #, Etc.)				13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE			14 CITY		15 STATE	16 ZIP CODE		
08 YEARS OF OPERATION		09 NAME OF OWNER									

**III. PREVIOUS OPERATOR (S)** *(List most recent first; provide if different from owner)***PREVIOUS OPERATOR'S PARENT COMPANIES** *(if applicable)*

01 NAME N/A			02 D & B NUMBER			10 NAME N/A			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)				04 SIC CODE		12 STREET ADDRESS (P.O. Box, RFD #, Etc.)				13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE			14 CITY		15 STATE	16 ZIP CODE		
08 YEARS OF OPERATION		09 NAME OF OWNER									

01 NAME			02 D & B NUMBER			10 NAME			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)				04 SIC CODE		12 STREET ADDRESS (P.O. Box, RFD #, Etc.)				13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE			14 CITY		15 STATE	16 ZIP CODE		
08 YEARS OF OPERATION		09 NAME OF OWNER									

01 NAME			02 D & B NUMBER			10 NAME			11 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)				04 SIC CODE		12 STREET ADDRESS (P.O. Box, RFD #, Etc.)				13 SIC CODE	
05 CITY		06 STATE	07 ZIP CODE			14 CITY		15 STATE	16 ZIP CODE		
08 YEARS OF OPERATION		09 NAME OF OWNER									

**IV. SOURCES OF INFORMATION** *(Cite specific references, e.g., state files, sample analysis, reports)*

NUS Corporation, FIT 3. Preliminary assessment. TDD No. F3-9008-01, June 27, 1991.

**EPA****POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 9 - GENERATOR/TRANSPORTER INFORMATION****I. IDENTIFICATION**01 STATE  
PA02 SITE NUMBER  
2898**II. ON-SITE GENERATOR***ORIGINAL*  
(HAC)

01 NAME N/A	02 D & B NUMBER	
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)	04 SIC CODE	
05 CITY	06 STATE	07 ZIP CODE

**III. OFF-SITE GENERATOR(S)**

01 NAME N/A	02 D & B NUMBER	01 NAME	02 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)	04 SIC CODE	03 STREET ADDRESS (P.O. Box, RFD #, Etc.)	04 SIC CODE		
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE
01 NAME N/A	02 D & B NUMBER	01 NAME	02 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)	04 SIC CODE	03 STREET ADDRESS (P.O. Box, RFD #, Etc.)	04 SIC CODE		
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE

**IV. TRANSPORTER(S)**

01 NAME N/A	02 D & B NUMBER	01 NAME	02 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)	04 SIC CODE	03 STREET ADDRESS (P.O. Box, RFD #, Etc.)	04 SIC CODE		
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE
01 NAME	02 D & B NUMBER	01 NAME	02 D & B NUMBER		
03 STREET ADDRESS (P.O. Box, RFD #, Etc.)	04 SIC CODE	03 STREET ADDRESS (P.O. Box, RFD #, Etc.)	04 SIC CODE		
05 CITY	06 STATE	07 ZIP CODE	05 CITY	06 STATE	07 ZIP CODE

**V. SOURCES OF INFORMATION** (Cite specific references, e.g., state files, sample analysis, reports)

NUS Corporation, FIT 3. Preliminary assessment. TDD No. F3-9008-01, June 27, 1991.



**EPA** ORIGINAL  
(Red)

**POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 10 - PAST RESPONSE ACTIVITIES**

**I. IDENTIFICATION**

01 STATE  
PA

02 SITE NUMBER  
2898

**H. PAST RESPONSE ACTIVITIES**

01 ☐ A. WATER SUPPLY CLOSED

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ B. TEMPORARY WATER SUPPLY PROVIDED

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ C. PERMANENT WATER SUPPLY PROVIDED

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ D. SPILLED MATERIAL REMOVED

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ E. CONTAMINATED SOIL REMOVED

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ F. WASTE REPACKAGED

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ G. WASTE DISPOSED ELSEWHERE

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ H. ON-SITE BURIAL

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ I. IN SITU CHEMICAL TREATMENT

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ J. IN SITU BIOLOGICAL TREATMENT

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ K. IN SITU PHYSICAL TREATMENT

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ L. ENCAPSULATION

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ M. EMERGENCY WASTE TREATMENT

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ N. CUTOFF WALLS

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ O. EMERGENCY DIKING/SURFACE WATER DIVERSION

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ P. CUTOFF TRENCHES/SUMP

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

01 ☐ Q. SUBSURFACE CUTOFF WALL

02 DATE \_\_\_\_\_

03 AGENCY \_\_\_\_\_

04 DESCRIPTION

N/A

**EPA**

**POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 10 - PAST RESPONSE ACTIVITIES**

**I. IDENTIFICATION**

01 STATE

PA

02 SITE NUMBER

2898

**II. PAST RESPONSE ACTIVITIES (Continued)**01 ☐ R. BARRIER WALLS CONSTRUCTED

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ S. CAPPING/COVERING

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ T. BULK TANKAGE REPAIRED

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ U. GROUT CURTAIN CONSTRUCTED

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ V. BOTTOM SEALED

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ W. GAS CONTROL

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ X. FIRE CONTROL

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ Y. LEACHATE TREATMENT

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ Z. AREA EVACUATED

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ 1. ACCESS TO SITE RESTRICTED

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ 2. POPULATION RELOCATED

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

01 ☐ 3. OTHER REMEDIAL ACTIVITIES

02 DATE

03 AGENCY

04 DESCRIPTION

N/A

**III. SOURCES OF INFORMATION** (Cite specific references, e.g., state files, sample analysis, reports)

NUS Corporation, FIT 3. Preliminary assessment. TDD No. F3-9008-01, June 27, 1991.

**EPA**ORIGINAL  
(Red)**POTENTIAL HAZARDOUS WASTE SITE  
SITE INSPECTION REPORT  
PART 11 - ENFORCEMENT INFORMATION****I. IDENTIFICATION**01 STATE  
PA02 SITE NUMBER  
2898**II. ENFORCEMENT INFORMATION**01 PAST REGULATORY ACTION ☒ YES ☐ NO

## 02 DESCRIPTION OF FEDERAL, STATE, LOCAL REGULATORY/ENFORCEMENT ACTION

On September 5, 1972, the facility was found to be in violation of the Clean Streams Law by the Pennsylvania Department of Environmental Resources after an inspection revealed several discharges to the Schuylkill River without a permit. These discharges included deionized wastewater, cooling water, sink wastewater, water softener waste, and septic tank seepage.

An October 31, 1973 PA DER inspection found the facility still in violation of the Clean Streams Law.

On February 8, 1974, the facility was fined \$250 by PA DER for violations of the Clean Streams Law.

On April 12, 1975, a PA DER inspection found that the facility's septic system was overflowing. Two discharges from the site were sampled and analyzed. Both contained unacceptable BOD levels and one contained ethyl glycol alcohol. This was a violation of the Clean Streams Law.

In a March 19, 1976 administrative conference with the facility, PA DER required the pumping of the existing septic tank as soon as the new tile field was in operation and the submission of a NPDES permit application.

On July 13, 1979, PA DER required the facility to submit a permit for the discharges to groundwater via a limestone pit.

On April 27, 1983, a PA DER inspection revealed the discharge of non-contact compressor cooling water to a storm sewer tributary to the Schuylkill River.

On December 6, 1984, PA DER conducted an inspection of the facility in response to citizens' complaints of eye and nasal irritation caused by the cooling water discharge. The inspection revealed that the facility was processing an ammonia-based cleaner. A sample of the discharge contained 5 ppb phenol and a BOD concentration greater than 405 ppm. On May 8, 1985, PA DER found the facility in violation of its NPDES permit when a sample of the non-contact cooling water discharge contained a BOD concentration of 11.5 ppm.

On August 13, 1985, PA DER found the facility to be in violation of the Clean Streams Law because of various potential sources of pollution at the site that could impact the NPDES discharge.

On March 14 and 27, 1991, EPA Emergency Response performed a removal site inspection after NUS Corporation, FIT 3 found numerous tanks and drums on site during a preliminary assessment on February 19, 1991. Many of the drums observed by the FIT (approximately 75; 25 contained material) had been moved into an on-site warehouse by the site owner. The large on-site holding tanks were found to be empty. Numerous drums and containers were observed inside the various buildings. No removal or enforcement action was performed. The OSC planned to work with the site owner to address safety concerns posed by the unrestricted site access and the materials haphazardly stored on site.

**III. SOURCES OF INFORMATION** (Cite specific references, e.g., state files, sample analysis, reports)

NUS Corporation, FIT 3. Preliminary assessment. TDD No. F3-9008-01, June 27, 1991.  
EPA file information.

## **SECTION 6.0**

## 6.0 REFERENCES FOR SECTIONS 1.0 THROUGH 5.0

1. United States Geological Survey. Phoenixville, Pennsylvania Quadrangle, 7.5 Minute Series. Topographic Map. 1955, photorevised 1983. Combined with Collegeville, Pennsylvania Quadrangle, 7.5 Minute Series. Topographic Map. 1966, photorevised 1983; Pottstown, Pennsylvania Quadrangle, 7.5 Minute Series. Topographic Map. 1956, photorevised 1968 and 1973; and Sassamansville, Pennsylvania Quadrangle, 7.5 Minute Series. Topographic Map. 1957, photorevised 1990.
2. HALLIBURTON NUS Environmental Corporation, FIT 3. Preliminary Assessment. TDD No. F3-9008-01, June 27, 1991.
3. NUS Corporation, FIT 3. Site discovery report; Township of Limerick. TDD No. F3-9003-06, June 22, 1990.
4. HALLIBURTON NUS Environmental Corporation. Site inspection; site visit. Project No. 3263-09, January 30, 1992.
5. Oestreicher, Eugene, Linfield Industrial Park, with (b) (4) NUS FIT 3. Telecon. January 23, 1991.
6. Herman, David L., Publicker Industries, Incorporated, to Joseph A. Feola, Pennsylvania Department of Environmental Resources. Correspondence. June 9, 1986.
7. Bretherick, Robert P., Publicker Packaging Services, Division of Publicker Industries, Incorporated, to Marilyn Shup, Pennsylvania Department of Environmental Resources. Correspondence. September 9, 1985.
8. Paly, John F., Pennsylvania Department of Environmental Resources, to H.E. Huber, Continental Distilling Corporation. Correspondence. September 5, 1972.
9. Laregrina, James, Pennsylvania Department of Environmental Resources, to Robert Bretherick, Publicker Industries, Incorporated. Correspondence. January 22, 1985.

ORIGINAL  
(Red)

Site Name: Linfield Industrial Park  
TDD No.: 3263-09

10. Pennsylvania Department of Environmental Resources, to Lee Vernon, Publicker Industries, Incorporated. Correspondence. January 22, 1974.
11. Simple, Andrew K., Pennsylvania Department of Environmental Resources, to Lee Vernon, Continental Distilling Corporation. Correspondence. February 27, 1974.
12. Minkle, Richard L., Pennsylvania Department of Environmental Resources, to Richard Grace, Continental Distilling Corporation. Correspondence. February 10, 1975.
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## **SECTION 7.0**

## 7.0 LABORATORY DATA

### 7.1 SAMPLE DATA SUMMARY

The attached data summary contains only analytes which were identified as detected in at least one sample. The complete list of compounds analyzed for, their results, and the associated detection limits are located as an appendix. Results for tentatively identified compounds appear following the organic data section of this report.

The following codes are used in the data summary to indicate the confidence in the laboratory results:

#### CODES RELATING TO IDENTIFICATION

(confidence concerning presence or absence of analytes):

- U = Not detected. The associated number indicates approximate sample concentration necessary to be detected.
- (NO CODE) = Confirmed identification.
- B = Not detected substantially above the level reported in laboratory or field blanks.
- R = Unreliable result. Analyte may or may not be present in the sample. Supporting data necessary to confirm result.
- N = Tentative identification. Consider present. Special methods may be needed to confirm its presence or absence in future sampling efforts.

#### CODES RELATED TO QUANTITATION

(can be used for both positive results and sample quantitation or detection limits):

- J = Analyte present. Reported value may not be accurate or precise.
- K = Analyte present. Reported value may be biased high. Actual value is expected to be lower.
- L = Analyte present. Reported value may be biased low. Actual value is expected to be higher.
- UJ = Not detected, quantitation or detection limit may be inaccurate or imprecise.
- UL = Not detected, quantitation or detection limit is probably higher.

#### OTHER CODES

- Q = No analytical result.

SITE NAME: Linfield Industrial Park  
TDD NUMBER: F3-3253-09  
LAB NAMES: ORG: Compu INORG: Keytx

SAMPLING DATE(S): 1/28/92 STATE/COUNTY CODE: PA-091  
CASE NUMBER: 17144 EPA NUMBER: 2898

[illegible][illegible]

Chemical Name	Concentration (ppm)	Mass (g)	Volume (L)
10.00 1,2,4-trichlorobenzene	1.00 J	100.00 J	100.00 J
10.00 naphthalene		180.00 J	
10.00 2-methyl naphthalene			
10.00 acenaphthylene			
10.00 acenaphthene			
10.00 4-nitrophenol			
10.00 dibenzofuran		120.00 J	
10.00 diethyl phthalate		180.00 B	
10.00 fluorene		270.00 J	
10.00 phenanthrene		2400.00	
10.00 anthracene		550.00 J	
10.00 di-n-butyl phthalate		150.00 J	
10.00 pyrene		72.00 J	
10.00 butyl benzyl phthalate		79.00 B	
10.00 benzolanthracene		360.00 J	
10.00 chrysene		320.00 J	
10.00 bis(2-ethylhexyl) phthalate		130.00 J	
10.00 benzobiphenylene		76.00 B	
10.00 benzofluoranthene		2300.00	
10.00 benzol(a)pyrene		210.00 J	
10.00 benzol(b)fluoranthene		1800.00	
10.00 benzol(k)fluoranthene		210.00 J	
10.00 benzol(a)anthracene		280.00 B	
10.00 benzol(b)anthracene		53.00 B	
10.00 benzol(k)anthracene		250.00 J	
10.00 benzol(a)pyrene		170.00 J	
10.00 indeno(1,2,3-cd)pyrene		1800.00	
10.00 dibenz(a,h)anthracene		1200.00	
10.00 benzo(g,h,i)perylene		410.00 J	
10.00 carbazole		86.00 J	
10.00 2-methyl naphthalene		95.00 J	
10.00 2-ethyl naphthalene		1200.00 N	
10.00 2-methyl anthracene		87.00 J	
10.00 2-methyl phenanthrene		91.00 J	

TYPE OF DATA	PESTICIDES	DILUTION FACTOR	1.0	1.000.0	1.000.0	1.000.0	1.0	46.0	51.0
1.0	1.000.0	1.000.0	1.000.0	1.000.0	1.000.0	1.000.0	1.0	46.0	51.0



SITE NAME: Linfield Industrial Park  
 TDO NUMBER: F1-3263-03  
 LAB NAMES: CRG: Compu INORG: Keytx

SELECTED SAMPLE ORDER  
 SAMPLING DATE(s): 1/26/92 STATE/COUNTY CODE: PA-091  
 CASE NUMBER: 17744 EPA NUMBER: 2896

SAMPLE NUMBER:	CHY02	CHY03	CHY04	CHY05	CHY06	CHY07	CHY08	CHY09	CHY10	CHY24	CHY11	CHY12	CHY13	
SAMPLE ID:	PSW-1	(b) (6)	(b) (6)	SW-1	SW-2	SW-3	SW-4	SW-5	SW-6	AQ-0	SD-1	SD-2	SD-3	
LOCATION:	CUHWC			UPSTREAM FRO	MIDSTREAM ON	ONSITE DRAIN	DOWNSTREAM	CUHWC; SCHUY	DUPLICATE OF	AQUEOUS	UPSTREAM ON	MID-STREAM	ON-SITE DRAIN	
	potable	potable	potable	N SCHUYLKILL	SCHUYLKILL R	AGE DITCH; 30	SAMPLE FROM	LKILL RIVER	SW-4 SAME	BLANK	SCHUYLKILL R	ON SCHUYL R.	DITCH; 30 FT	
	no treat	no treat	no treat	RIVER, UNDER	AT FORMER	FEET FROM CO	SCHUYLKILL	INTAKE, ACRO	LOCATION AS		UNDER MAIN	AT FORMER	FROM CONFL.	
	clear, no odr	clear, no odr	clear, no odr	MAIN ST. BRIG	NPDES DISCHG	NFL. W/ SCHU. R	RIVER	SS FROM SITE	SW-4		ST. BRIDGE	NPDES DISCHG	W/ SCHUYL R.	
PH:				6.9	6.7	6.7	6.9				6.3	6.4	6.5	
FIELD MEASUREMENTS:	none>bkg hnu	none>bkg hnu	none>bkg hnu	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	
PERCENT SOLIDS:											71.0%	45.0%	64.0%	
TYPE OF DATA: ***** PESTICIDES														
DILUTION FACTOR:	1.0	1.0	1000.0	1000.0	1000.0	1.0	1000.0	1000.0	1000.0	1.0	45.0	73.0	51.0	
DET. LIMIT	SAMPLE NUMBER:	CHY02	CHY03	CHY04	CHY05	CHY06	CHY07	CHY08	CHY09	CHY10	CHY24	CHY11	CHY12	CHY13
CRQL (μg/L)	UNITS:	ug/l	ug/l	ng/l	ng/l	ng/l	ug/l	ng/l	ng/l	ng/l	ug/l	ug/kg	ug/kg	ug/kg
0.05 alpha-BHC												3.40 R	0.38 R	
0.05 beta-BHC														
0.05 delta-BHC												1.50DR		
0.05 gamma-BHC (lindane)												34.00 R		
0.05 heptachlor													0.13 B	0.14 B
0.05 aldrin												4.40 R	1.10 R	0.33 R
0.05 heptachlor epoxide												1.70DR		0.24 R
0.05 endosulfan I														
0.10 dieldrin														0.46 R
0.10 4,4'-DDE													4.40 P	
0.10 endrin												3.90DR		1.60 R
0.10 endosulfan II													2.50 R	0.32 R
0.10 4,4'-DDD					5.30 R	7.10 R		4.60 R	5.00 R	4.90 R		2.80 R	4.80 R	
0.10 endosulfan sulfate													4.70 R	
0.10 4,4'-DDT													2.20 R	
0.50 methoxychlor								14.00 R				7.80 R	15.00 R	2.70 B
0.10 endrin ketone												3.80 R		0.65 R
0.05 alpha-chlordane				2.90 R								5.30 R	4.10 R	0.26 R
0.05 gamma-chlordane												2.70DR	3.50 R	0.13 R
1.00 brochlor-1260														
1.00 endrin aldehyde													2.60 P	

Comments: \*\*\*\*\*

sample quant. limit=dil. factor x CRQL D= result reported from diluted re-analysis  
 C= confirmed by GC/MS

SITE NAME: Linfield Industrial Park  
TDD NUMBER: F3-3283-09  
LAB NAMES: ORG: Compu INORG: Keytx

SELECTED SAMPLE ORDER  
SAMPLING DATE(s): 1/28/92 STATE/COUNTY CODE: PA-081  
CASE NUMBER: 17744 EPA NUMBER: 2898

SAMPLE NUMBER:		CHY14	CHY15	CHY16	CHY17	CHY18	CHY19	CHY22	CHY20	CHY21	CHY23	CHY26	CHY25
SAMPLE ID:		SD-4	S-1	S-2	S-3	S-4	S-5	S-6	S-6	S-7	S-9	S-10	S-0
LOCATION:		DOWNSTREAM SAMPLE ON SCHUYLKILL RIVER	AT BASE OF TRANSFORMER ON NORTH SID E OF DISTIL.	AT DRUMS OUTSIDE OF SHED WEST OF DISTILLERY.	INSIDE ABOVE GROUND TANK FARM ON EAST SIDE BLDG.2	AT CONFLU. W/2 CONCRETE DRAIN. TROUGH HS SW WAREHS	AT CORRUGATED PIPE OUTFALL SOUTH AND WE ST OF WAREHS	DUPLICATE OF S-5 SAME LOCATIO N AS S-5	SE EDGE OF TRASH DUMP WEST OF WAREHOUSES	10-FT. SOUTH OF DRUMS BY TRASHDUMP WE ST OF WAREHS	IN SOUTHERN AREA OF SITE AN AREA OF STRESS. VEG.	ACROSS MAIN STREET FROM SITE	TRIP BLANK
PH:		6.2	6.7	6.0	6.6	6.1	6.2	6.4	6.7	6.6	6.6	6.0	
FIELD MEASUREMENTS:		NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE
PERCENT SOLIDS:		44.0%	81.0%	38.0%	78.0%	72.0%	76.0%	80.0%	65.0%	65.0%	78.0%	68.0%	
TYPE OF DATA: ***** VOLATILES													
DILUTION FACTOR:		2.3	1.2	2.6	1.3	1.4	1.3	1.2	1.5	1.5	1.3	1.5	1.0

DET. LIMIT	SAMPLE NUMBER:	CHY14	CHY15	CHY16	CHY17	CHY18	CHY19	CHY22	CHY20	CHY21	CHY23	CHY26	CHY25
CPQL (=IDL)	UNITS:	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/l
10.00	methylene chloride	54.00B	49.00B	120.00B	45.00B	52.00B	47.00B	56.00B	70.00B	51.00B	68.00B	83.00B	150.00
10.00	acetone	39.00B	61.00B	420.00B	28.00B	25.00B	57.00B	61.00B	69.00B	86.00B	37.00B	73.00B	
10.00	2-butanol			160.00									
10.00	1,1,1-trichloroethane												
10.00	4-methyl-2-pentanone			340.00									
10.00	toluene			120.00									
10.00	ethylbenzene			100.00									
10.00	total xylenes			690.00									
TYPE OF DATA: ***** SEMIVOLATILES													
DILUTION FACTOR:		74.0	1000.0	1700.0	84.0	45.0	86.0	41.0	750.0	250.0	42.0	49.0	9.0

DET. LIMIT	SAMPLE NUMBER:	CHY14	CHY15	CHY16	CHY17	CHY18	CHY19	CHY22	CHY20	CHY21	CHY23	CHY26	CHY25
CPQL (=IDL)	UNITS:	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/l
10.00	phenol												
10.00	4-methylphenol												
10.00	1,2,4-trichlorobenzene		2300.00 J										
10.00	naphthalene			4500.00 J			150.00 J	71.00 J	2500.00 J				
10.00	2-methylnaphthalene			8000.00 J			95.00 J	75.00 J	790.00 J				
10.00	acenaphthylene												
10.00	acenaphthene				99.00 J	89.00 J	440.00 J	150.00 J	9000.00	350.00 J			
50.00	4-nitrophenol												
10.00	dibenzofuran						130.00 J	61.00 J	2900.00 J				
10.00	diethyl phthalate										54.00 B		
10.00	fluorene			2600.00 J		66.00 J	180.00 J	89.00 J	4100.00 J				
10.00	phenanthrene	330.00 J	2900.00 J	5500.00 J	620.00 J	1400.00 J	2300.00 J	900.00 J	52000.00 J	3500.00 J		77.00 J	
10.00	anthracene	77.00 J				140.00 J	450.00 J	150.00 J	14000.00	920.00 J			
10.00	di-n-butyl phthalate												
10.00	fluoranthene	790.00	7100.00 J	4900.00 J	1200.00 J	3900.00 J	4100.00 J	1500.00 J	72000.00 J	8600.00 J	79.00 J	170.00 J	
10.00	pyrene	450.00 J	4900.00 J	11000.00 J	920.00 J	3300.00 J	2000.00 J	1300.00 J	54000.00 J	3500.00 J	63.00 J	99.00 J	
10.00	butylbenzyl phthalate	130.00 B								470.00 B			
10.00	benzofluoranthene	420.00 J	3000.00 J	5200.00 J	620.00 J	1700.00	1400.00	730.00 J	33000.00	4500.00		72.00 J	
10.00	chrysene	340.00 J	3800.00 J	4300.00 J	610.00 J	2100.00	1500.00	880.00 J	32000.00	3800.00	44.00 J	96.00 J	
10.00	bis(2-ethylhexyl)phthalate	350.00 B	1300.00 B	7500.00 B					1600.00 B				
15.00	benzo(b)fluoranthene	360.00 J	7600.00 J	3900.00 J	1500.00 J	4500.00 J	2100.00 J	1600.00 J	46000.00 J	6600.00 J	64.00 J	105.00 J	
10.00	benzo(k)fluoranthene	240.00 J			1500.00 J		830.00 J	570.00 J	15000.00 J	3800.00 J		63.00 J	
10.00	benzofluorene	360.00 J	3700.00 J	3700.00 J	480.00 J	1900.00 J	1500.00 J	550.00 J	30000.00 J	6900.00 J		86.00 J	
10.00	indeno(1,2,3-cd)pyrene	220.00 J	2400.00 J		500.00 J	1700.00	1100.00	370.00 J	15000.00	3800.00			
10.00	dibenz(a,h)anthracene	79.00 J			200.00 J	350.00 J	400.00 J	180.00 J	5300.00 J	1700.00 J			
10.00	benzo(g,h,i)perylene	190.00 J	2100.00 J	7500.00 J	520.00 J	1800.00	1000.00	910.00 J	7600.00	3500.00			
10.00	carbazole				100.00 J	180.00 J	330.00 J	120.00 J	11000.00	710.00 J			
TYPE OF DATA: ***** PESTICIDES													
DILUTION FACTOR:		76.0	910.0	86.0	42.0	46.0	43.0	41.0	56.0	51.0	42.0	48.0	0.0

DET. LIMIT	SAMPLE NUMBER:	CHY14	CHY15	CHY16	CHY17	CHY18	CHY19	CHY22	CHY20	CHY21	CHY23	CHY26	CHY25
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SITE NAME: Linfield Industrial Park  
TDD NUMBER: F3-3263-03  
LAB NAMES: ORG: Compu INORG: Keytx

SELECTED SAMPLE ORDER  
SAMPLING DATE(s): 1/28/92 STATE/COUNTY CODE: PA-091  
CASE NUMBER: 17744 EPA NUMBER: 2898

SAMPLE NUMBER:	CHY14	CHY15	CHY16	CHY17	CHY18	CHY19	CHY22	CHY20	CHY21	CHY23	CHY26	CHY25
SAMPLE ID:	SD-4	S-1	S-2	S-3	S-4	S-5	S-8	S-6	S-7	S-9	S-10	S-0
LOCATION:	DOWNSTREAM SAMPLE ON SCHUYLKILL RIVER	AT BASE OF TRANSFORMER ON NORTH STD E OF DISTIL.	AT DRUMS OUTSIDE OF SHED WEST OF DISTILLERY.	INSIDE ABOVE GROUND TANK FARM ON EAST SIDE BLDG.2	AT CONFLU. W/2 CONCRETE DRAIN. TROUGH HS SW WAREHS	AT CORRUGATD PIPE OUTFALL SOUTH AND WE ST OF WAREHS	DUPLICATE OF S-5 SAME LOCATIO N AS S-5	SE EDGE OF TRASH DUMP WEST OF WAREHOUSES	10-FT. SOUTH OF DRUMS BY TRASHDUMP WE ST OF WAREHS	IN SOUTHERN AREA OF SITE AN AREA OF STRESS. VEG.	ACROSS MAIN STREET FROM SITE	TRIP BLANK
PH:	6.2	6.7	6.0	6.6	6.1	6.2	6.4	6.7	6.6	6.6	6.0	
FIELD MEASUREMENTS:	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE
PERCENT SOLIDS:	44.0%	81.0%	38.0%	78.0%	72.0%	76.0%	80.0%	65.0%	65.0%	78.0%	68.0%	
TYPE OF DATA: *****	PESTICIDES *****											
DILUTION FACTOR:	76.0	810.0	86.0	42.0	46.0	43.0	41.0	50.0	51.0	42.0	48.0	

DET. LIMIT	SAMPLE NUMBER:	CHY14	CHY15	CHY16	CHY17	CHY18	CHY19	CHY22	CHY20	CHY21	CHY23	CHY26	CHY25
CRQL (*=IDL)	UNITS:	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	n/a
0.05 alpha-BHC		0.36 R	11.00 R	1.10 R			0.08 R						
0.05 beta-BHC				3.90 R					1.30 R		0.87 R	0.28 R	
0.05 delta-BHC		5.50DR											
0.05 gamma-BHC (lindane)		34.00DR	3.40 R	4.40 R	14.00 R					0.18 R		0.09 R	
0.05 heptachlor				0.16 B	16.00 R		0.15 B		1.80 R	0.81 B	0.23 B	0.19 B	
0.05 aldrin		4.30 R	12.00 R	2.60 R	14.00 R	0.36 R	0.39 R		5.60 R	1.40 R		0.66 R	
0.05 heptachlor epoxide		5.60 R			0.19 R	0.54 R	0.41 R		2.30 R				
0.05 endosulfan I		14.00DR		0.37 R									
0.10 dieldrin		3.30DR		37.00 R	31.00 R	0.27 R	0.10 R	6.20 R	14.00 R	2.40 R		0.26 R	
0.10 4,4'-DDE				13.00 R		2.80 R	1.10 R	1.60 R	11.00 R	6.40 R		0.80 R	
0.10 endrin		4.50DR		46.00 R	39.00 R	2.60 R	1.50 R		34.00 R		1.40 R	3.30 R	
0.10 endosulfan II				64.00 R	2.20 R	2.60 R	0.48 R		20.00 R		0.41 R	0.28 R	
0.10 4,4'-DDD		4.30 R											
0.10 endosulfan sulfate		0.69 B		19.00 R									
0.10 4,4'-DDT		1.80DR			31.00 R	0.48 R		2.30 R		13.00 R		0.92 R	
0.50 methoxychlor		10.00 R		180.00 R	31.00 R	18.00 R	6.50 R	26.00 R	120.00 R	110.00 R	1.40 B		
0.10 endrin ketone		4.60 R						0.66 R					
0.05 alpha-chlordane		6.20 R		11.00 R				3.50 R		2.40 R			
0.05 gamma-chlordane		3.30 R		11.00 P	0.09 R	0.23 R		2.10 R	2.80 R	5.70 R			
1.00 aroclor-1260			300000.00DC							460.00			
1.00 endrin aldehyde		3.10 R		53.00 R					17.00 R	20.00 R	0.68 R		

Comments: \*\*\*\*\*  
sample quant.limit=dil.factor X CRQL    D=result reported from diluted re-analys  
C= confirmed by GC/MS

SITE NAME: Linfield Industrial Park  
 TDD NUMBER: F3-3263-09  
 LAB NAMES: ORG: Compu INORG: Keytz

SELECTED SAMPLE ORDER  
 SAMPLING DATE(s): 1/28/92 STATE/COUNTY CODE: PA-091  
 CASE NUMBER: 17744 EPA NUMBER: 2898

SAMPLE NUMBER:	MCJE14	MCJE15	MCJE16	MCJE17	MCJE18	MCJE19	MCJE22	MCJE20	MCJE21	MCJE23	MCJE26	
SAMPLE ID:	S0-4	S-1	S-2	S-3	S-4	S-5	S-8	S-6	S-7	S-9	S-10	
LOCATION:	DOWNSTREAM SAMPLE IN SCHUYLKILL RIVER	AT BASE OF TRANSFORMER ON NORTH SID E OF DISTIL.	AT DRUMS OUTSIDE OF SHED WEST OF DISTILLERY	INSIDE ABOVE GROUND TANK FARM ON EAST SIDE BLDG.2	AT CONFLU. OF 2 CONCRETE DRAIN. TROUGH HS; SW WAREHS	AT CORRUG. PIPE OUTFALL SOUTH AND WE ST OF WAREHS	DUPLICATE OF S-5 SAME LOCATIO N AS S-5	SE EDGE OF TRASH DUMP WEST OF WAREHOUSES	10-FEET SOUT H OF DRUMS NEAR TRASHOW W. OF WAREHSE	IN SOUTHERN PORTION OF SITE; AREA OF STRESS. VEG.	ACROSS MAIN STREET FROM SITE	
PH:	6.2	6.7	6.0	6.6	6.1	6.2	6.4	6.7	6.6	6.6	6.0	
FIELD MEASUREMENTS:	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	NONE	
PERCENT SOLIDS:	43.2%	63.6%	43.7%	61.8%	76.0%	83.1%	74.8%	64.7%	74.9%	77.4%	67.9%	
TYPE OF DATA: ***** INORGANICS *****												
DILUTION FACTOR: : GFAA	0.463	0.239	0.458	0.244	0.263	0.241	0.268	0.309	0.267	0.258	0.295	
: ICP	0.463	0.239	0.458	0.244	0.263	0.241	0.268	0.309	0.267	0.258	0.295	
: Hg	1.160	0.600	1.140	0.611	0.658	0.602	0.668	0.773	0.668	0.646	0.736	
: CN	0.120	0.060	0.114	0.061	0.066	0.060	0.067	0.077	0.066	0.065	0.074	
DET. LIMIT	SAMPLE NUMBER:	MCJE14	MCJE15	MCJE16	MCJE17	MCJE18	MCJE19	MCJE22	MCJE20	MCJE21	MCJE23	MCJE26
CRQL (≡IDL)	UNITS:	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
14.00 aluminum		10500.00	6810.00	2140.00	4970.00	3400.00	4170.00	5470.00	3780.00	5090.00	8230.00	6790.00
3.00 antimony										11.00		
2.00 arsenic		3.80 B	372.00	62.40 J	3.90 B	6.70 B	5.40 B	10.40	9.00	33.30	5.60 B	3.40 B
1200.00 barium		154.00	158.00	407.00	60.60	52.70	66.00	113.00	80.90	268.00	111.00	44.60
1.00 beryllium		2.10	0.70	0.55	0.71	0.51	0.58	0.78	0.53	0.65	1.00	0.72
3.00 cadmium		2.20	2.60	2.80					1.10	2.30		
7000.00 calcium		4560.00	1270.00	2220.00	1520.00	2110.00	1820.00	2670.00	12400.00	17600.00	1730.00	382.00
3.00 chromium		43.20	29.10	63.90	12.40	10.60	8.10	8.20	38.60	45.10	15.20	14.20
4.00 cobalt		59.70	14.30	9.60	8.60	6.40	7.50	8.70	8.80	14.80	11.10	9.30 B
3.00 copper		105.00	42.70 J	71.50 J	15.40 J	10.00 J	7.10 J	17.00 J	81.70 J	388.00 J	20.10 J	13.50 J
6.00 iron		23000.00	37700.00	39000.00	12800.00	10800.00	9880.00	11600.00	10300.00	67800.00	20900.00	18100.00
1.00 lead (anal. by GFAA)		100.00	349.00	4810.00	35.70	18.20	16.90	22.20 J	253.00	358.00	20.20 J	17.80 J
37.00 magnesium		3100.00	1830.00	784.00	1040.00	937.00	1140.00	1070.00	4750.00	4860.00	1890.00	1250.00
2.00 manganese		1540.00	411.00	464.00	811.00	318.00	351.00	469.00	437.00	801.00	1090.00	366.00
0.20 mercury		0.35	0.18	2.90					0.23	0.33		
22.00 nickel		75.20	26.90	10.60	9.80	7.40	7.10	12.70	10.70	56.50	14.40	10.70
722.00 potassium		970.00	822.00		782.00	397.00	483.00	483.00	534.00	1100.00	883.00	582.00
2.00 selenium					0.71 B	0.63 B	0.60 B	0.59 B	0.98 B			0.82 B
30.00 sodium		195.00	129.00	52.40	64.50	43.50	115.00	91.00	63.40	219.00	41.10	32.10
4.00 vanadium		28.20	41.10	26.50	15.70	19.60	12.30	19.80	12.70	73.10	31.50	27.20
2.00 zinc		497.00	1230.00	651.00	200.00	169.00	201.00	286.00	226.00	1150.00	61.80	129.00
10.00 cyanide			0.68	1.80								

Comments: \*\*\*\*\*

D= result reported from diluted re-analysis  
 sample det. limit = dil. factor x IDL

C= confirmed by GC/MS

SITE NAME: Linfield Industrial Park  
 TOD NUMBER: F3-3263-09  
 LAB NAMES: ORG: Compu INORG: Keytx

SELECTED SAMPLE ORDER  
 SAMPLING DATE(s): 1/28/92 STATE/COUNTY CODE: PA-091  
 CASE NUMBER: 17744 EPA NUMBER: 2898

SAMPLE NUMBER: MCJE02 MCJE03 MCJE04 MCJE05 MCJE06 MCJE07 MCJE08 MCJE10 MCJE09 MCJE24 MCJE11 MCJE12 MCJE13														
SAMPLE ID: PSM-1 HW-1 HW-2 SW-1 SW-2 SW-3 SW-4 SW-6 SW-5 AQ-0 SD-1 SD-2 SD-3														
LOCATION: CUHWC Zwolak Brown, UPSTREAM MIDSTREAM ON ONSITE DRAIN DOWN STREAM DUPLICATE CUHWC: SCHUYL AQUEOUS UPSTREAM ON MID-STREAM ON-SITE DITC														
potable no treat no treat no treat FROM SCHUYL SCHUYLKILL R AGE DITCH: 30 SAMPLE FROM SCHUYLKILL OF SW-4 KILL RIVER BLANK SCHUYLKILL R ON SCHUYL R. H 30-FT. FRO														
clear, no odr clear, no odr clear, no odr ER MAIN ST. NPDES DISCHG NFLU. W/ SCH. R PI-VER SAME LOCATIO INTAKE. ACR- OSS FROM SIT ST. BRIDGE NPDES DISC. SCHUYLKILL R														
PH: 6.9 6.7 6.7 5.9 6.8 6.3 6.4 6.5														
FIELD MEASUREMENTS: none>bkg hnu none>bkg hnu none>bkg hnu NONE NONE NONE NONE NONE NONE 51.2% 40NE 50.4% 60.0%														
PERCENT SOLIDS:														
TYPE OF DATA: ***** INORGANICS *****														
DILUTION FACTOR: : GFAA 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 0.391 0.397 0.333														
: ICP 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 0.391 0.397 0.333														
: Mg 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 0.977 0.992 0.833														
: CN 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 1.000 0.098 0.099 0.083														
DET. LIMIT	SAMPLE NUMBER:	MCJE02	MCJE03	MCJE04	MCJE05	MCJE06	MCJE07	MCJE08	MCJE10	MCJE09	MCJE24	MCJE11	MCJE12	MCJE13
CRQ1 (x=IDL)	UNITS:	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l	ug/l	mg/kg	mg/kg	mg/kg
14.00 aluminum		22.90 B	15.10 B	32.50 B	134.00 J	937.00 J	1090.00 J	1560.00 J	266.00 J	119.00 J	15.19	6080.00	8190.00	12400.00
3.00 antimony														
2.00 arsenic		10.80 B	8.10 B	5.80 B	6.80 B	7.20 B	7.20 B	8.90 B	6.40 B	6.70 B	5.40	7.00 B	6.10 B	7.80 B
120.00 barium		95.50	248.00	254.00	33.00 J	55.60 J	65.30 J	93.90 J	33.80 J	30.30 J		120.00	118.00	146.00
1.00 beryllium								1.10				1.10	1.30	2.20
3.00 cadmium												1.40	1.40	
7000.00 calcium		68700.00	38700.00	48400.00	38000.00	38900.00	14500.00	44100.00	37900.00	33100.00	29.10	10200.00	4070.00	1800.00
3.00 chromium						3.70 J		7.10 J				31.60	31.50	19.30
4.00 cobalt						6.70 J		29.00 J	6.40 J			38.80	32.90	31.00
3.00 copper		5.00	188.00	68.00	5.50 J	23.80 J	5.00 J	55.60 J	7.40 J	4.50 J		74.20	58.30	61.00
6.00 iron		9.40 B	52.50 B	21.90 B	202.00 J	1380.00 J	930.00 J	3470.00 J	547.00 J	132.00 J	12.50	14900.00	17800.00	17500.00
1.00 lead (anal. by GFAA)						19.00 J	1.10 J	59.90 J	2.70 J			113.00 J	62.70	33.30 J
37.00 magnesium		15100.00	15300.00	16800.00	12400.00	12700.00	5790.00	14300.00	12400.00	10800.00	42.60	5420.00	3110.00	2580.00
2.00 manganese				2.70	123.00 J	507.00 J	51.70 J	1210.00 J	379.00 J	113.00 J		1340.00	1230.00	291.00
0.20 mercury												0.23	0.30	0.50
22.00 nickel								32.60				40.30	40.40	55.20
22.00 potassium		774.00	1000.00	1300.00	3050.00 J	3300.00 J	2270.00 J	3770.00 J	2620.00 J	2780.00 J		629.00	560.00	869.00
2.00 selenium		2.00 B			2.00 B	2.00 B			3.30 B	2.50 B		1.50 B		0.96 B
33.00 sodium		14700.00	21900.00	17750.00	23300.00	22600.00	6150.00	24100.00	22300.00	18800.00		203.00	221.00	80.60
4.00 vanadium								6.40				23.30	22.80	23.20
2.00 zinc		14.40 B	63.30 B	39.50 B	48.50 B	176.00 J	27.70 B	301.00 J	61.40 B	76.00 B	15.80	384.00	288.00	239.00
10.00 cyanide						11.70								

Comments: \*\*\*\*\*

0= result reported from diluted re-analysis  
 C= confirmed by GC/MS  
 sample det. limit= dil. factor x IDL

## **7.2 ORGANIC DATA VALIDATION REPORT: CASE 17744**

### **7.2.1 Overview**

Case 17744 consisted of 14 low-level solid samples and 10 low-level aqueous samples submitted to Compuchem Laboratories (COMPU) for volatiles, semivolatiles, and pesticide/polychlorinated biphenyl (PCB) analyses. One field blank and two duplicate pairs were included in this sample set. One trip blank was analyzed for volatile compounds only. The samples were analyzed as a Contract Laboratory Program (CLP) Routine Analytical Service (RAS).

### **7.2.2 Summary**

These samples were analyzed according to the March 1990 Statement of Work (SOW), and the data were fully reviewed using the National Functional Guidelines for Organic Data Review, Multi-Media, Multi-Concentration, February 1988, with Regional Modifications. (Areas examined in detail are included in the Support Documentation appendix to this report.) All samples were successfully analyzed for all target compounds. Detection capability for most analytes was demonstrated by meeting criteria for spike and surrogate recoveries and instrument tuning and calibration. The laboratory performed all analyses in accordance with CLP RAS protocols.

### **7.2.3 Major Issues**

Issues affecting or changing the confidence concerning the presence or absence of analytes and issues related to severe quantitative bias are discussed below.

- All results for methylene chloride, acetone, diethyl phthalate, di-n-butyl phthalate, butylbenzyl phthalate, and bis(2-ethylhexyl) phthalate (DEHP) have been flagged as artifacts of blank contamination (B), except where these compounds were detected in the field blanks, samples CHY24 and CHY25. The results for these common laboratory contaminants were not substantially higher in the samples than in all associated blanks.

- Several single-peak pesticide results have also been flagged (B) because concentrations were within five times the concentrations for corresponding peaks detected in blank analyses. The affected compounds and associated results are as follows:

Compound	Affected Results
heptachlor	CHY12, CHY13, CHY16, CHY19, CHY21, CHY23, and CHY26
endosulfan sulfate	CHY14
methoxychlor	CHY13 and CHY23

- Many pesticide compounds were reported, but, because none were found at levels high enough to allow GC/MS confirmation, the confidence concerning the presence of pesticides depends upon the retention time of each reported compound being unique (i.e., different from any unrelated detectable compound that might happen to be in the sample). Unfortunately, there were so many closely spaced peaks of similar sizes in the chromatograms of solid samples, there is a high probability that unrelated compounds could have eluted at retention times similar to the target pesticides and thereby produced false positives for the reported pesticide compounds. Furthermore, in many cases, there were severe disagreements between the concentrations of a pesticide calculated from injection of the same extract on two different GC columns, which suggests that the larger of the two peaks must be attributable to interferences.

Therefore, it cannot be determined if any individual target pesticide compound reported in these samples is actually present, versus an artifact attributable to unrelated organic compounds that are not part of the CLP target compound list. All target pesticide results have been flagged (R), unreliable, to indicate that these compounds may or may not be present. Even groups of related pesticide compounds reported together in samples CHY11, CHY12, CHY13, and CHY14 are not considered reliable. [For example, the components of technical-grade chlordane (alpha chlordane, gamma chlordane, heptachlor, and heptachlor epoxide) were reported together, but because there were continuous closely spaced peaks adjacent to these target compounds that were of similar size, technical chlordane constituents were not visible above the random interference pattern that was superimposed over the target compound retention time zones. In addition, several of the identifications of chlordane and DDT analogs exhibited greater than four-fold discrepancies between the concentrations detected on two GC columns.] The Support Documentation attachments contain example chromatograms and comparisons of retention times and concentrations detected on the two GC columns. Further information would be useful to determine whether any of the reported pesticides are actually present.

- The reviewer corrected the quantitation of results for benzo(b)fluoranthene and benzo(k)fluoranthene in all samples where both of these isomers were reported, except CHY17 and CHY22. In these samples, the laboratory identified and reported these isomers as a co-eluting pair and assumed that equal amounts of both isomers were present. The laboratory personnel did not attempt a manual integration of the separate chromatographic peaks in order to obtain more accurate results for the individual isomers. In some cases, the mass-selective chromatogram indicated that only the benzo(b) fluoranthene isomer was present above the laboratory's reporting limit of 1 ng. In samples CHY11, CHY13, CHY15, CHY16, CHY18, and CHY23, the reviewer deleted the results for benzo(k)fluoranthene from the data summary. In samples CHY12, CHY14, CHY19, CHY20, CHY21, and CHY26, the reviewer changed the percentage of area attributed to each isomer. Because the reviewer's recalculations for the areas associated with each isomer are estimates, all results changed by the reviewer should be considered estimated. The Support Documentation appendix contains all chromatograms, quantitation lists for the samples and associated standards, and re-calculations.
- Very low recoveries (less than 10 percent) were reported for the tribromophenol surrogate in sample CHY22. This sample was re-analyzed and exhibited similar recoveries. Detection capability for acid-fraction compounds, in particular multihalogenated compounds such as pentachlorophenol, is considered unreliable for this sample. Because quality control data were similar for base-neutral surrogates, the reviewer reported the higher of the initial and re-analysis results for each base-neutral compound that was not affected by blank contamination and the lower of the two results for compounds within the range affected by blank contamination.
- Very low matrix spike/matrix spike duplicate (MS/MSD) recoveries (two and 10 percent, respectively) were obtained for pentachlorophenol in solid sample CHY18. Corresponding tribromophenol surrogate recoveries in these analyses were 12 and 53 percent, respectively. Therefore, detection capability for pentachlorophenol is considered unreliable in all solid samples.



- The result for benzo(g,h,i)perylene has been flagged as tentative (N) in sample CHY11. The mass spectrum for this compound did not contain all the characteristic ions that are present in the reference standard. The base peak appeared to be obscured by interferences, and the lower intensity half-mass ions were not visible. Because many other polynuclear aromatic compounds are present in this sample, and because this compound was confidently identified in nearly every other sample at this site, this compound should be considered present at this location. Further information may be useful in confirming the presence of this compound at this sampling location.

#### **7.2.4 Minor Issues**

Data qualifications related to precision, accuracy, and bias in reported results and quantitation limits are discussed below.

- The recoveries for the acid semivolatile surrogate compounds 2-fluorophenol, tribromophenol, and 2-chlorophenol were low in solid sample CHY16. (The recovery for 2-fluorophenol was below contractual limits.) Detection limits for acid-fraction compounds are considered biased low in this sample.
- The laboratory diluted the semivolatile fractions of samples CHY18 and CHY20 because levels for some target compounds exceeded the calibration range of the instrument. The reviewer reported the highest levels for all compounds that were within the calibration range of the instrument that were not considered artifacts of blank contamination (such as phthalates). For compounds that exhibited instrument levels that were below the calibration range, the results that were closer to the lower limit of the calibration range (10 ng) were reported. For compounds present at levels affected by blank contamination, the lowest of the two results was reported. Note that diethyl phthalate was not reported in sample CHY18 because it was detected only in the diluted analysis.
- The recovery for the tribromophenol surrogate was low for the initial, undiluted analysis of sample CHY18. (This recovery was acceptable for the diluted re-analysis.) The detection limits for multihalogenated phenols (especially pentachlorophenol) are considered biased low for the initial analysis of this sample. However, the elevated detection limits reported from the diluted analysis are considered valid, with the exception of pentachlorophenol, which was previously qualified unreliable.

ORIGINAL  
(100)

Site Name: Linfield Industrial Park  
Project No.: 3263-09

- Low recoveries for the tribromophenol surrogate were obtained for samples CHY17, CHY19, and CHY20. Detection limits for multihalogenated phenols (especially pentachlorophenol) are considered biased low for these samples.
- Low recoveries were noted for pyrene for the MS/MSD analysis of sample CHY18. Because the recoveries for the other base-neutral compounds were acceptable, and the terphenyl surrogate recoveries were acceptable, it is possible that sample inhomogeneity may have contributed to these low recoveries. A direction of bias cannot be firmly established. All solid results for pyrene have been flagged as estimated (J) and may be either imprecise or biased slightly low.
- The field duplicate samples CHY19 and CHY22 exhibited imprecision in results above the quantitation limit for phenanthrene, fluoranthene, and benzo(a)pyrene. Results for these compounds are considered estimated and have been flagged (J) in all surface soil samples (CHY15 through CHY23 and CHY26).
- A high percent difference between the average initial calibration relative response factor (RRF) and the continuing calibration RRF for benzo(g,h,i)perylene was observed on the day that sample CHY16 was analyzed. The result for this compound should be considered estimated in this sample.
- The laboratory diluted and re-analyzed the pesticide/PCB fraction of samples CHY11, CHY14, and CHY15 because of target compounds that exceeded the calibration range of the instrument. Based on an evaluation of both analyses for each sample, the reviewer reported the highest levels for all compounds that were detected in either analysis that were not rejected because of blank contamination. For sample CHY15, the results for Aroclor 1260 were reported from the diluted analysis because the instrument level result was within the range of the calibration standards for this PCB.
- The extraction for the semivolatile analysis of solid sample CHY16 occurred 19 days after sampling. Detection limits may be biased low for compounds that were not detected, although the degree of bias may be very minor for compounds known to be persistent in the environment [polycyclic aromatic hydrocarbons (PAHs), in particular]. The result for DEHP in this sample was previously flagged due to blank contamination. All other detected compounds were PAHs and were previously flagged (J), estimated, because levels are below the quantitation limit.

#### 7.2.5 Notes

- The laboratory extracts of the pesticide fraction of samples CHY15 and CHY21 were accidentally switched in the initial analyses, as revealed when high levels of PCBs were confirmed by GC/MS in the semivolatile extracts of the opposite sample. A re-extraction of these two samples was performed that confirmed that sample CHY15 contained high levels of PCBs and sample CHY21 contained only a very low level.
- The laboratory did not provide the second page of the semivolatile result form (Ic) for sample CHY02. Raw data indicated which results the laboratory intended to report, so the reviewer has reported the appropriate values on the data summary.
- Although flagged (B), the result for acetone may be real in sample CHY16. The instrument level result was higher than for any other sample, and the presence of 2-butanone (a related ketone that is often present when acetone is present in a sample) may also suggest that acetone may be indigenous to this sampling location. Further information may be useful in verifying if acetone is actually present at this location.
- The reviewer noted the presence of 4-nitrophenol, an unusual compound, in sample CHY07. This compound was confirmed by acceptable spectrum and retention time matching. Although this sample was also chosen for MS/MSD analysis, there was no evidence of any matrix spike compounds (other than 4-nitrophenol) in the unspiked analysis of CHY07.
- The results reported for Aroclor 1260 in samples CHY15 and CHY21 exhibited acceptable peak-pattern matching quality and retention time match for characteristic peaks on both chromatographic columns. In addition, penta- through heptachlorobiphenyls were also identified as tentatively identified compounds (TICs) in sample CHY21; these congeners are the major constituents of Aroclor 1260.
- Because the laboratory did not supply mass-selective chromatograms for sample CHY17, it cannot be determined whether only one compound or both benzo(b)fluoranthene and benzo(k)fluoranthene are present and in what relative quantities each isomer may be present.

(b) (4)

(b) (4)

Case 17744

## SAMPLE DATA SUMMARY: ORGANIC TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NUMBER	ANALYSIS FRACTION (VOA/BNA)	ESTIMATED CONCENTRATION		QUALIFIER CODE	COMPOUND NAME
		VALUE	UNITS		
					* There were no VOA TICS in any sample except CHY16 *
CHY16	VOA	1670	ug/kg	TOT/ISO	C <sub>3</sub> -alkyl benzene, C <sub>9</sub> H <sub>12</sub> (7)
↓	↓	21	↓	ISO	C <sub>4</sub> -alkyl benzene, C <sub>10</sub> H <sub>14</sub>
↓	↓	120	↓	ISO	C <sub>9</sub> H <sub>10</sub> , methyl styrene
CHY02	BNA	5	ug/l		aliphatic alcohol
"	"	8		TOT	sat'd HC (3)
CHY03	BNA				ND
CHY04	BNA				ND
CHY05	BNA	3	ug/l	UNK	pass. Nitrogen subst. aromatic
CHY06	BNA				ND
CHY07	BNA				ND
CHY08	BNA				ND
CHY09	BNA				ND
CHY10	BNA				ND

## DEFINITIONS OF QUALIFIER CODES:

**SUS = SUSPECTED FALSE POSITIVE RESULT:** Compound is either a common laboratory contaminant, or else a possible reaction byproduct (artifact) attributable to the chemical reagents used for sample preparation and analysis. This result is suspect even though this compound was not found in any associated blanks.

**UNK = UNKNOWN COMPOUND:** Library search result unreasonable or of very low matching quality.

**TOT = TOTAL CONCENTRATION REPORTED:** Represents the sum of several compounds detected all belonging to the same chemical class.

**ISO = OR ISOMER:** Compound identification is not selective for this isomer only. This result may instead represent the presence of a similar compound comprised of the same atoms bonded together in a different arrangement or substitution pattern.

**SAT HC = SATURATED HYDROCARBON**

**UNSAT HC = UNSATURATED HYDROCARBON**

**HC = HYDROCARBON**

**PAH = polynuclear aromatic hydrocarbon**

**SUB = SUBSTITUTED**

**MIX = MIXTURE OF 2 OR MORE COELUTANTS**

**ND = NONE DETECTED**

**SAMPLE DATA SUMMARY: ORGANIC TENTATIVELY IDENTIFIED COMPOUNDS**

see 17744

SAMPLE NUMBER	ANALYSIS FRACTION (VOA/BNA)	ESTIMATED CONCENTRATION		QUALIFIER CODE	COMPOUND NAME
		VALUE	UNITS		
CHY11	BNA	650	ug/kg	ISO	C <sub>15</sub> H <sub>10</sub> PAH, such as cyclopenta[1,2,3-cd]pyrene
		3000		TOT/ISO	C <sub>17</sub> H <sub>12</sub> PAH, such as methyl pyrene (3)
		920		ISO	C <sub>19</sub> H <sub>14</sub> PAH, such as methyl chrysene
		2300		ISO	C <sub>20</sub> H <sub>12</sub> PAH, such as benzo(e)pyrene
↓	↓	4100	↓	TOT	Sat'd HC (3)
CHY12	BNA	360	ug/kg	ISO	Methylbenzene sulfenamide - C <sub>10</sub> H <sub>9</sub> N <sub>2</sub> S <sub>2</sub> <sup>USP 70553</sup>
		1600		TOT	Carboxylic acid (3)
		5040		TOT	Sat'd HC (6)
		220		UNK	Unsat'd HC of unknown subst.
		730			aromatic amide
		660			aldehyde, such as octadecanal
↓	↓	2200	↓	UNK	possible sterol derivative
CHY13	BNA	460	ug/kg	TOT	aliphatic alcohols - (2)
		670		TOT	Carboxylic acids (2)
		520			Sat'd HC
↓	↓	1300	↓	TOT/UNK	Unknowns (2)
CHY14	BNA	150	ug/kg		Benzaldehyde
		450			Carboxylic acid
		2000		TOT/UNK	unsat'd HC of unknown subst. (3)
		4200		TOT	Sat'd HC (3)
		13,000			aliphatic alcohol
		600		ISO	C <sub>20</sub> H <sub>12</sub> PAH - benzo(e)pyrene
		9000		TOT/UNK	Unknowns (4)
↓	↓	3400	↓	UNK	possible sterol

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- ISO = OR ISOMER:** Compound identification is not selective for this isomer only. This result may instead represent the presence of a similar compound comprised of the same atoms bonded together in a different arrangement or substitution pattern.
- SAT HC = SATURATED HYDROCARBON**
- UNSAT HC = UNSATURATED HYDROCARBON**
- HC = HYDROCARBON**
- PAH = polynuclear aromatic hydrocarbon**
- SUB = SUBSTITUTED**
- MIX = MIXTURE OF 2 OR MORE COELUTANTS**
- ND = NONE DETECTED**

Case 17744

## SAMPLE DATA SUMMARY: ORGANIC TENTATIVELY IDENTIFIED COMPOUNDS

SAMPLE NUMBER	ANALYSIS FRACTION (VOA/BNA)	ESTIMATED CONCENTRATION		QUALIFIER CODE	COMPOUND NAME
		VALUE	UNITS		
CHY15	BNA	44,000	ug/kg	TOT	pentachlorophenyl (2)
		315,000		TOT	hexachlorobiphenyl (5)
		200,000		TOT	heptachlorobiphenyl (5)
		23,000		ISO	1,2,3,4,5,6-hexachloro-1,2,3,4,5,6-hexachlorobiphenyl
		43,000		TOT/ISO	decahydro dimethyl octyl naphthalene (2)
		130,000	ug/kg	TOT/UNK	Unknowns (4)
✓	✓	50,000	"	UNK	Unknown aromatic HC
CHY16	BNA	1.9 x 10 <sup>6</sup>	ug/kg	TOT	SAT'D HC (19)
"	"	82,000	"		cyclic HC - C <sub>17</sub> H <sub>34</sub> (alkylcyclohexane)
CHY17	BNA	1100	ug/kg	ISO	benzothiazole, C <sub>7</sub> H <sub>5</sub> N <sub>2</sub> S, 95-16-9
		510		ISO	Methylthiobenzothiazole (C <sub>7</sub> H <sub>5</sub> N <sub>2</sub> S, 615-22-5)
		1740		TOT	mercaptobenzothiazole, C <sub>7</sub> H <sub>5</sub> N <sub>2</sub> S, 149-30-4 (2)
		340		UNK	Nitrogen/sulfur subst. aromatic HC
		420		UNK	poss. Nitrogen subst. aromatic HC
		850		ISO	C <sub>20</sub> H <sub>12</sub> PAH, such as benzo(e)pyrene
✓	✓	4200		TOT/UNK	Unknowns (7)
CHY18	BNA	370	ug/kg		carboxylic acid
		180		ISO	cyclopentaphenanthrene
		270		ISO	anthracenodione
		2500		ISO	C <sub>20</sub> H <sub>12</sub> PAH, such as benzo(e)pyrene
✓	✓	18,000		TOT	Sat'd HC (3)

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**HC** = HYDROCARBON

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**SUB** = SUBSTITUTED

**MIX** = MIXTURE OF 2 OR MORE COELUTANTS

**ND** = NONE DETECTED

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# SAMPLE DATA SUMMARY: ORGANIC TENTATIVELY IDENTIFIED COMPOUNDS

Case 17744

SAMPLE NUMBER	ANALYSIS FRACTION (VOA/BNAL)	ESTIMATED CONCENTRATION		QUALIFIER CODE	COMPOUND NAME
		VALUE	UNITS		
CHY19	BNAL	350	ug/kg	ISO	cyclopenta phenanthrene
		350	ug	ISO	anthracene dione
		1300		ISO	C <sub>20</sub> H <sub>12</sub> PAH, such as benzo(e)pyrene
		4000	↓	TOT/UNK	UNKNOWN'S (7)
CHY20	BNAL	23,000	ug/kg	ISO	C <sub>10</sub> H <sub>6</sub> O <sub>3</sub> - hydroxynaphthalene dione
		12,000	↓	ISO	C <sub>15</sub> H <sub>10</sub> - cyclopenta phenanthrene
		4500		ISO	C <sub>15</sub> H <sub>12</sub> PAH - such as methylanthracene
		7600		ISO	C <sub>16</sub> H <sub>10</sub> O - benzonaphthofuran
		45,000		TOT/ISO	C <sub>17</sub> H <sub>12</sub> PAH - such as methyl pyrene (4)
		16,000		TOT/ISO	C <sub>17</sub> H <sub>14</sub> PAH - such as phenylmethyl naphthalene (2)
		18,000		ISO/TOT	C <sub>17</sub> H <sub>10</sub> O - benzoanthracene one (2)
		12000		ISO	C <sub>16</sub> H <sub>10</sub> O <sub>5</sub> - benzo naphthothioxophene
		25,000		TOT/ISO	C <sub>18</sub> H <sub>12</sub> PAH - such as benzo(c)phenanthrene (2)
		9800		ISO	C <sub>16</sub> H <sub>11</sub> N - benzo carbazole
		23,000		TOT/ISO	C <sub>19</sub> H <sub>14</sub> PAH - such as methyl chrysene (2)
		28,100		TOT/ISO	C <sub>20</sub> H <sub>12</sub> PAH - such as benzo(e)pyrene (2)
		17,000		UNK	mix of PAHs; m.w. 240, 246
		7600			aliphatic alcohol
		14,000	↓	TOT/UNK	UNKNOWN'S (2)
CHY21	BNAL	500	ug/kg	ISO	C <sub>15</sub> H <sub>10</sub> - cyclopenta phenanthrene
		6300	↓	TOT/ISO	C <sub>20</sub> H <sub>12</sub> PAH - such as benzo(e)pyrene (2)
		17000	↓	TOT	UNKNOWN'S (17)

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### **7.3 INORGANIC DATA VALIDATION REPORT: CASE NUMBER 17744**

#### **7.3.1 Overview**

The set of samples for Case 17744 contained 10 unfiltered aqueous samples and 14 solid samples that were analyzed through the CLP RAS by Keystone Engineering (KEYTX). Included in the sample set were an unfiltered field blank, an unfiltered field duplicate pair, and a solid field duplicate pair.

#### **7.3.2 Summary**

All analytes were successfully analyzed in all samples. Areas of concern with respect to data usability are listed according to the seriousness of the problem. These include the following items:

#### **7.3.3 Major Issues**

Issues affecting or changing the confidence concerning the presence or absence of analytes and issues related to severe quantitative bias are discussed below.

- The calibration blanks contained calcium and selenium at levels equal to or greater than the instrument detection limit (IDL). The solid preparation blank contained selenium at a level above the IDL, and the aqueous field blank contained aluminum, arsenic, calcium, iron, magnesium, and zinc at levels above the IDLs. The reported results for these analytes in the affected samples that are less than five times the highest blank concentration have been flagged (B) and may be artifacts of blank contamination (i.e., false positives).
- Results for selenium in samples MCJE02 and MCJE18 and lead in samples MCJE12 through MCJE15, MCJE17, and MCJE18 have been corrected by the reviewer. The laboratory miscalculated or misreported the affected results.

#### 7.3.4 Minor Issues

Data qualifications related to precision, accuracy, and bias in reported results and detection limits are discussed below.

- The inductively coupled plasma (ICP) serial dilution results for aluminum and iron exceeded the 10 percent control limit in aqueous sample MCJE07. Therefore, all reported results for these elements in similar samples have been qualified as estimated (J), except the field blank.
- A low correlation coefficient for arsenic occurred for the method of standard addition (MSA) analysis of one sample. The result is considered estimated (J) because a low correlation coefficient suggests a lack of linearity for the analytical curve generated by this method.
- Significant imprecision was noted between furnace and plasma results for lead in five solid samples. Because the laboratory satisfactorily performed all the required quality assurance/quality control (QA/QC) for the plasma analysis of lead, two valid sets of data are present for all solid samples. The lead results in the affected samples have been qualified as estimated (J); sample inhomogeneity may be responsible for the observed differences. Good agreement was noted between the two methods for the remainder of the solid samples and for aqueous sample MCJE08.
- Field duplicate imprecision for aluminum, barium, cobalt, copper, iron, lead, manganese, potassium, and zinc was noted between surface water samples MCJE08 and MCJE10. Therefore, results for these analytes should be considered estimated in all similar samples and have been qualified as (J). It is possible that a higher level of suspended particulate matter was present in the aliquot taken for digestion of sample MCJE08 than in the aliquot taken for sample MCJE10; this could explain why higher levels occurred for most analytes in MCJE08.
- Field duplicate imprecision also occurred for copper between solid samples MCJE19 and MCJE22. Therefore, results for this analyte should be considered estimated (J) in all similar surface soil samples.

### 7.3.5 Notes

- Good agreement between plasma and furnace was noted for arsenic in sample MCJE15. The instrument level was above 1,500 ug/l.
- Post-digestion spike (PDS) recoveries were low for selenium analyses in samples MCJE17, MCJE20, and MCJE22. This would suggest that the results for the affected samples may be biased low; however, all three results have been previously qualified because of blank contamination.
- The data were reviewed in accordance with the National Functional Guidelines for Evaluating Inorganic Analyses, with Region III Modifications.

Report prepared by (b) (4)  
(215) 971-0900

(b) (4)

Report reviewed by (b) (4)  
(215) 971-0900

(b) (4)

**TABLE 1A  
 SUMMARY OF QUALIFIERS ON DATA SUMMARY  
 AFTER DATA VALIDATION**

ANALYTE	SAMPLES AFFECTED	POSITIVE VALUES	BIAS	COMMENTS*
Al	MCJ302, MCJE03, and MCJE04	B	high	(A) (15.1 ug/l)
As	MCJE02 through MCJE14, MCJE17 through MCJE19, MCJE23, and MCJE26	B	high	(A) (5.4 ug/l)
Fe	MCJE02 through MCJE04	B	high	(A) (12.5 ug/l)
Se	all positive results	B	high	(A) (2.7 ug/l)
Zn	MCJE02 through MCJE05, MCJE07, MCJE09, and MCJE10	B	high	(A) (15.8 ug/l)
Al	MCJE05 through MCJE10	J	---	(B) (24.5 percent)
Fe	MCJE05 through MCJE10	J	---	(B) (17.5 percent)
As	MCJE16	J	---	(C) (0.7442)
Al	MCJE05 through MCJE10	J		(D) (142 percent)
Ba	MCJE05 through MCJE10	J		(D) (94 percent)
Co	MCJE05 through MCJE10	J		(D) (128 percent)
Cu	MCJE05 through MCJE10	J		(D) (153 percent)
Fe	MCJE05 through MCJE10	J		(D) (146 percent)
Pb	MCJE05 through MCJE10	J		(D) (183 percent)
Cu	MCJE15 through MCJE23 and MCJE26	J		(D) (82 percent)
Mn	MCJE05 through MCJE10	J		(D) (105 percent)
K	MCJE05 through MCJE10	J		(D) (36 percent)
Zn	MCJE05 through MCJE10	J		(D) (132 percent)
Pb	MCJE11, MCJE13, MCJE22, MCJE23, and MCJE26	J		(E) (68, 74, 99, 66, and 113 percent)

\*Comments are defined in table 1B.

**TABLE 1B**  
**CODES USED IN COMMENTS COLUMN**

- A = The blanks contained a result for this analyte that was greater than the IDL. (The highest blank value is in parentheses.) The reported results were less than five times this blank value and may be biased high.
- B = The ICP serial dilution result was outside the control limits. (The percent difference between the initial and diluted result is in parentheses.) The reported results are estimated.
- C = Because of a low correlation coefficient (the coefficient is in parentheses), the positive result be should be considered estimated.
- D = Field duplicate imprecision occurred for this analyte (the relative percent difference is in parentheses). The reported results are considered estimated.
- E = Imprecision between the graphite furnace and plasma values occurred for lead in this sample (the relative percent differences between these results for each sample are in parentheses, respectively). The result is considered estimated.

## **SECTION 8.0**

## 8.0 TOXICOLOGICAL EVALUATION

### 8.1 SUMMARY

Metals were detected at notable concentrations in soil from the transformer area, the drums west of the distillery, and the trash dump area. It could be prudent to avoid contact with the high-level lead and arsenic concentrations because neither noncarcinogenic nor carcinogenic risks can be ruled out. Polychlorinated biphenyls (PCBs) were detected above recommended guidelines in the soil sample taken from the base of the transformer; mainly because of carcinogenic risk, it could be prudent to avoid contact with this soil. It may be prudent to avoid prolonged contact with the soil from the southeastern edge of the trash dump because of notable polycyclic aromatic hydrocarbon (PAH) concentrations, although adverse effects are not necessarily indicated. PAHs can also pose some increased cancer risk. Reported concentrations of 1,2,4-trichlorobenzene (1,2,4-TCB), toluene, ethylbenzene, xylenes, ketones, cyanide, cadmium, and antimony in soil are not expected to pose significant hazards.

Cadmium, phenol, 1,1,1-trichloroethane (1,1,1-TCEA), 4-methylphenol (4-MP), and PAHs were detected in river sediment at concentrations not expected to pose significant noncarcinogenic direct-contact hazards. The potential for bioconcentration of sediment cadmium by aquatic organisms could not be ruled out.

4-Nitrophenol (4-NP) was detected in drainage ditch water at a concentration not expected to pose a significant environmental or human health threat. Copper, iron, lead, zinc, and cyanide were detected in the river above Ambient Water Quality Criteria (AWQCs); potential effects on sensitive aquatic organisms cannot be ruled out. Metals in the sample taken from the river at a drinking water intake were below health-based criteria and are not expected to pose a significant hazard. Metals in the river water are not expected to pose a significant noncarcinogenic health threat to swimmers. A theoretical increase in overall cancer risk due to lead and beryllium in river water cannot be ruled out.

Xylenes were detected in the public supply well far below drinking water criteria; no adverse effects are indicated. Sodium was detected in one home well above a concentration recommended to minimize sodium intake; adverse effects on the general population are not indicated.



## 8.2 SUPPORT DOCUMENTATION

### 8.2.1 Organics

PCBs were confidently and tentatively detected in soil from the base of the transformer. The confidently detected result was reported at 300,000 ug/kg. PCBs were also detected in soil south of the trash dump area at 460 ug/kg. PCBs were at one time used in transformers because they are stable, thermoplastic, and nonflammable.<sup>1</sup> PCBs are also extremely persistent and stable in the environment and can bioconcentrate in the food chain.<sup>1</sup> EPA has recommended guidelines of 25,000 ug/kg for industrial or restricted access areas and 10,000 ug/kg for residential or unlimited access areas.<sup>2</sup> Recently, a protective quantifiable concentration of 1,000 ug/kg has been recommended for residential areas.<sup>3</sup> The reported concentration at the on-site "hot spot" exceeds all recommended guidelines.

Prolonged contact with high concentrations of PCBs can be associated with chloracne and liver damage.<sup>1</sup> However, nonoccupational human toxicity due to PCBs has not been reported for people in the United States.<sup>4</sup> The main concern for human exposure to PCBs is the classification of these compounds as Group B2 (probable human) carcinogens. For a 15-kilogram child incidentally ingesting 200 mg of this soil 350 days per year for six years (an extremely conservative scenario), the estimated excess upper-bound cancer risk would be approximately  $2\text{E-}3$ .<sup>5</sup> For a 70-kilogram adult (100 mg soil per day for 30 years), the risk would be approximately  $1\text{E-}3$ .<sup>5</sup> Because any contact with carcinogens can theoretically increase overall cancer risk, it would be prudent to avoid contact with this soil.

PAHs were confidently and tentatively identified in soil and sediment throughout the site, as well as in background soil. PAHs are common environmental contaminants because they occur in coal, tar, and the products of the combustion of organic material.<sup>6</sup> Coal tar, which is rich in PAHs, was once used for the extraction of dye compounds such as benzene, toluene, naphthalene, phenol, and carbazole.<sup>7</sup> These compounds, as well as other compounds associated with dye synthesis, were found on and near the site. Dye synthesis is not known to have occurred on site. The plastics industry is reported to be an important market for dyes; plastics manufacturing may have occurred on site (see section 1.3).<sup>7</sup> PAHs can also be naturally occurring up to as much as approximately 10,000 ug/kg.<sup>8</sup> At this site, PAHs were confidently detected up to 62,400 ug/kg in on-site soil, with a much higher concentration of 406,190 ug/kg at sample S-6 (southeastern edge of trash dump), compared to 758 ug/kg in background soil. Sample S-6 was described as containing large amounts of organic material (see sample log).

PAHs have been associated with skin irritation at high concentrations.<sup>1</sup> Although such effects are not necessarily indicated from reported concentrations, it may be prudent to avoid prolonged contact with the highest-concentration soil. Even if a 15-kilogram child were exposed daily to 200 mg of the most contaminated soil or sediment, significant noncarcinogenic effects would not be indicated due to these PAHs, separately or together, based on a comparison between estimated intakes and available risk reference doses (RfDs) [total Hazard Quotient (HQ) (ratio of intake to RfD), 0.07].<sup>5,9</sup> Estimated total cancer risks from PAHs in on-site soil ingested at this rate would be up to approximately 4E-4, using worst-case (maximum) concentrations for each PAH.<sup>5,9,10</sup> This can be compared to the cancer risk of 6E-7, which could be estimated for background soil.<sup>5,9,10</sup> The maximum estimated oral cancer risk for sediment using the same exposure scenario would be approximately 2E-5.

An examination of the pattern of sediment contamination reveals that concentrations of total PAHs were lowest in the drainage ditch and downstream samples and highest in the upstream sample. In spite of their lipophilicity, PAHs tend to be metabolized by many aquatic organisms, reducing bioconcentration potential and secondary exposure to fish consumers, if any.<sup>11</sup>

1,2,4-TCB was detected in one on-site soil sample at 2,300 ug/kg. The 1,2,4-TCB, which can be used as a dielectric fluid in transformers, was detected in the sample obtained at the base of the transformer.<sup>1</sup> 1,2,4-TCB can cause skin irritation, neurological effects, and liver damage at high (such as occupational industrial) exposures.<sup>1</sup> Comparing a conservative estimate of 1,2,4-TCB intake from exposure to this soil (15-kilogram child ingesting 200 mg soil 350 days per year for six years) to the RfD, it can be seen that adverse noncarcinogenic effects would not be anticipated (HQ = 0.003).<sup>9</sup>

Toluene (120 ug/kg), ethylbenzene (100 ug/kg), and xylenes (690 ug/kg) were detected in one soil sample (S-2, at the drums west of the distillery). These compounds are volatile solvents that can be irritants and neurotoxicants at high concentrations.<sup>1</sup> They tend to be associated with fuels or paints and are fairly mobile in the environment.<sup>1</sup> Xylenes and ethylbenzene can also be used in the manufacture of plastics, and toluene can be a component of dyeing wastes.<sup>12</sup> Comparison between conservatively estimated intakes and the RfDs reveals that noncarcinogenic impacts are not indicated for these compounds, separately or together (HQs for toluene, ethylbenzene, and xylenes are 8E-6, 1E-5, and 4E-6, respectively).<sup>5</sup>

Two ketones were also detected in the soil sample taken from the drum area west of the distillery: 2-butanone (160 ug/kg) and 4-methyl-2-pentanone (340 ug/kg). These ketones are solvents that can have irritant effects.<sup>1</sup> However, comparison of conservatively estimated intakes to the RfDs indicates that these compounds would not be expected to pose a significant noncarcinogenic hazard, separately or together (HQs of 2-butanone and 4-methyl-2-pentanone are 4E-5 and 8E-5, respectively).<sup>5,9</sup>

Phenol (75 ug/kg) and 4-MP (95 ug/kg) were detected in midstream river sediment. Phenol is a strong irritant that can be used as a disinfectant at low concentrations.<sup>1</sup> It is a potential component of dyeing wastes.<sup>12</sup> 4-MP is a disinfectant that can be used in the manufacture of chemicals, dyes, and plastics.<sup>1</sup> 4-MP is also an irritant.<sup>1</sup> Comparison of conservatively estimated intakes to the RfDs indicates that these compounds would not be expected to pose a significant noncarcinogenic direct contact hazard, individually or together (HQs of phenol and 4-MP are 1.6E-6 and 2.4E-5, respectively).<sup>9</sup> These compounds are generally not very persistent in the aquatic environment, and biodegradation is usually significant.<sup>11,13</sup>

1,1,1-TCEA was detected in upstream river sediment at 3 ug/kg. This solvent can affect the skin, cardiovascular system, and nervous system at high levels.<sup>1</sup> However, at the reported trace concentration, no significant impacts are indicated.<sup>5,9</sup> This compound is not very persistent in the aquatic environment.<sup>11,13</sup>

4-NP was detected in drainage ditch surface water at approximately 1 ug/l, well below the chronic fresh-water Lowest Observed Effects Level (LOEL) of 150 ug/l for nitrophenols.<sup>14</sup> Nitrophenols are used in pesticide and dye synthesis.<sup>1</sup> 4-NP is reported to affect aquatic organisms at concentrations exceeding 4,000 ug/l.<sup>13</sup> The 4-NP concentration is also well below the level that would be of concern for human ingestion of the water, using an RfD of 6.2E-2 mg/kg/day and assuming one liter per day consumption by a 15-kilogram child (HQ = 7E-5).<sup>15</sup>

Xylenes were detected in the PSW-1 sample at approximately 1 ug/l. The properties of xylenes were discussed previously. The PSW-1 concentration is well below the enforceable Maximum Contaminant Level (MCL) and purely health-based MCL Goal (MCLG) of 10,000 ug/l.<sup>16</sup> No adverse effects are indicated for this trace concentration of xylenes.

An aliphatic alcohol (approximately 5 ug/l) and saturated hydrocarbons (approximately 8 ug/l) were also tentatively identified in PSW-1. The alcohols are generally solvents that can be irritants at high concentrations.<sup>17</sup> The saturated hydrocarbons could include solvents, plant constituents, etc. No specific toxicological assessment of unknown compounds can be made.

Methylbenzene sulfonamide, a fungicide, was tentatively identified in the midstream sediment sample. Few data are available for this compound. However, an oral wild bird LD50 (dose lethal to 50 percent of an experimental population) of 75 mg/kg has been reported.<sup>18</sup> In soil sample S-3 (taken from the tank farm), other sulfur-containing compounds such as benzothiazole and nitrogen-containing compounds were tentatively identified. Benzothiazole has an oral mouse LD50 of 900 mg/kg.<sup>18</sup> Sulfur- and nitrogen-containing compounds are widely used in dye synthesis.<sup>7,19</sup>

### 8.2.2 Inorganics

Metals were detected at notable concentrations in four soil samples: S-1 (at the base of the transformer), S-2 (at the drums outside the shed west of the distillery), S-6 (at the southeastern edge of the trash dump), and S-7 (south of the drums near the trash dump).

Lead was detected up to 358 mg/kg in soil samples (excluding S-2). Lead has been known to affect the hematopoietic, renal, and nervous systems and is classified as a Group B2 carcinogen.<sup>9,18</sup> Children, because of their high absorption rate and developing nervous systems, are especially vulnerable.<sup>6</sup> These soil lead concentrations are below a recommended clean-up guideline of 500 to 1,000 mg/kg.<sup>20</sup> Lead binds strongly to soil, decreasing its availability. Significant impacts are not necessarily expected for these concentrations of lead in soil, although it is generally desirable to minimize all lead exposure. Lead was detected at 4,810 mg/kg in soil sample S-2. Inadvertent ingestion of 200 mg soil would result in a lead intake of 962 ug. Daily lead intake from food, air, and dust is reported to range from 60.2 (rural nonsmoker) to 118.2 (urban smoker) ug/day for adults and from 46.6 (rural) to 137.6 (urban) ug/day for children.<sup>21</sup> Chronic exposure could be undesirable, as increases in blood lead have been reported from chronic lead exposure (35 days or longer).<sup>22</sup> Effects from increased blood lead range from enzyme inhibition through anemia and encephalopathy in extreme cases.<sup>6</sup> It would be prudent to avoid prolonged contact with the apparent "hot spot" at S-2, especially for children.

Arsenic was detected in on-site soil up to 372 mg/kg; the highest concentration was detected at sample S-1. Arsenic has been used in pesticides, pharmaceuticals manufacturing, glass, textiles, and metal alloys.<sup>1</sup> Arsenic is noted for skin toxicity and gastrointestinal and neurological symptoms.<sup>1</sup> The concentrations in soil samples S-1, S-2, and S-7 would result in an HQ greater than 1.0, assuming frequent direct contact (incidental ingestion of 200 mg soil by a 15-kilogram child).<sup>9</sup> Therefore, it would be prudent to avoid direct contact with soils from these areas, although arsenic is generally strongly bound to soil and the scenario is extremely conservative in assuming 100 percent absorption. Arsenic is also classified as a Group A (human) oral carcinogen.<sup>5</sup> Using the oral cancer slope factor (CSF) for arsenic of 1.75 per mg/kg/day and the above conservative scenario (also assuming exposure duration of six years and the very conservative exposure frequency of 350 days per year), an upper-bound cancer risk of approximately  $7\text{E-}4$  can be estimated for children if they were in almost daily contact with the highest-level soil sample.<sup>5</sup> For adults (same scenario, but assuming 100 mg soil per day, 70-kilogram body weight, and 30 years exposure duration), the HQ for the highest-level arsenic only would exceed 1.0; the estimated cancer risk would be approximately  $4\text{E-}4$ . Because of the high CSF for arsenic, it would be prudent to avoid contact with the highest-level soils and is generally desirable to minimize contact with arsenic. However, arsenic, like lead, is a naturally occurring element for which it is virtually impossible to attain zero exposure.

Cyanide was detected in soil at the base of the transformer and at the drums west of the distillery up to 1.8 mg/kg. Cyanides can be used in electroplating and other industries.<sup>1</sup> At high concentrations (e.g., occupational), cyanides can cause nausea, skin irritation, and eventually interfere with the ability of red blood cells to release oxygen to tissues (its fatal mechanism).<sup>1</sup> However, cyanide is not generally considered to be a major environmental problem because of its lack of persistence; it is readily metabolized at low concentrations.<sup>23</sup> Based on the conservative estimated intake for children, an oral HQ of  $1\text{E-}3$  can be calculated.<sup>5,9</sup> Significant noncarcinogenic impacts due to cyanide are not expected for this type of contact with soil.

Antimony was detected at a notable concentration (11 mg/kg) in S-7. Antimony is used in metal alloys, plasticizers, paints and glazes, and ammunition and explosives.<sup>1</sup> It can cause skin irritation and gastrointestinal effects at high concentrations.<sup>1</sup> Using the conservative scenario for child soil exposure, the reported concentration of antimony is not expected to pose a significant noncarcinogenic hazard ( $\text{HQ} = 0.3$ ).<sup>5</sup>

Cadmium was detected up to 2.6 mg/kg in on-site soil and up to 2.2 mg/kg in sediment from the Schuylkill River (cadmium was detected upstream, midstream, and downstream of the site, but was not detected in the on-site drainage ditch). Cadmium is a common environmental contaminant that can cause toxicity to the kidneys, prostate, and blood after high-level industrial exposure.<sup>1</sup> The reported cadmium concentrations would not be expected to pose a significant noncarcinogenic hazard (HQ up to 0.07, using the conservative child exposure scenario).<sup>5</sup> Cadmium tends to bind strongly to sediment and was not detected in the surface water. Bioconcentration of cadmium can be significant for certain freshwater species, including caddisflies, brook trout, and mosquitofish [reported bioconcentration factors (BCFs) range from 2 to 4,190].<sup>24</sup> The reported cadmium concentrations in fish and shellfish used as food in a typical diet range from 0.1 to 1 mg/kg.<sup>6</sup>

Notable concentrations of metals were not reported for soil sample S-9, where stressed vegetation was observed.

The close proximity of state game lands introduces the possibility that ecological receptors may be exposed to on-site metals. Potential receptors include deer, opossum, squirrels, rabbits, and mice. Most of these animals would be expected to venture only rarely onto the most contaminated areas of the site, because of the lack of cover from trees and the presence of houses and roads in the northern area. Because of the proximity of the river, animals such as ducks, geese, and gulls may be present; however, these birds are expected to spend more time in and very near the water. Both lead and arsenic have been reported to affect rodents, large mammals such as cattle, and fowl at sufficient concentrations.<sup>25</sup> In fowl, the effects of arsenic toxicity include inflammation of the gizzard and liver, and the effects of lead toxicity include gastrointestinal effects and neuropathy.<sup>25</sup> *Gastrointestinal effects and neurotoxicity are reported in larger animals for both metals.*<sup>25</sup> With the information available at this stage of site investigations, it is difficult (and beyond the current scope) to estimate effects on such animals or their consumers, but a slight possibility of exposure to these animals cannot be ruled out.

Metals were also detected at notable concentrations in surface water samples obtained from the Schuylkill River and the drainage ditch. Aluminum was detected in all samples, ranging from 119 ug/l (at the CUHWC intake) to up to 1,560 ug/l (downstream Schuylkill River), in excess of the AWQC of 87 ug/l.<sup>26</sup> For other metals, the only exceedances of AWQCs occurred in the midstream and downstream river samples: copper (23.8 ug/l midstream, 55.6 ug/l downstream; AWQC: 12 ug/l), iron (1,380 ug/l midstream, 3,470 ug/l downstream; AWQC: 1,000 ug/l), lead (19 ug/l midstream, 59.9 ug/l downstream; AWQC: 3.2 ug/l), zinc (176 ug/l midstream, 301 ug/l downstream; AWQC: 110 ug/l), and cyanide (11.7 ug/l midstream; AWQC: 5.2 ug/l).<sup>14</sup> Where AWQCs are exceeded, potential effects on sensitive aquatic species cannot be ruled out. Lead is especially undesirable in an aquatic environment because of its tendency to bioconcentrate (reported BCFs range from approximately 40 in vertebrate fish to 1,700 in a snail).<sup>27</sup> Potential effects on fish consumers, if any, cannot be specifically assessed without fish-tissue analysis.

Because the river is ultimately used for drinking water purposes at the CUHWC intake, it is possible to examine this sample (SW-5) for pretreatment metals. In this sample, the only metals that exceeded drinking water criteria were aluminum (119 ug/l) and manganese (113 ug/l), which exceeded the non-health-based aesthetic Secondary MCLs (SMCLs) of 50 ug/l.<sup>28</sup> The metals in this sample would not be expected to produce adverse health effects on the general population, even if the water were consumed in this condition.

It is also possible to examine the surface water results with regard to recreational use, such as swimming. The most conservative approach includes comparison of the samples with drinking water criteria or RfDs. SMCLs were not considered, but the health-based RfDs for chemicals such as aluminum, iron, and manganese were considered. Even if the river water were consumed at the rate of one liter per day by a 15-kilogram child, neither the individual HQs nor the total Hazard Index (HI) (sum of HQs) would exceed 1.0.<sup>5,9</sup> In actuality, the estimated intake of water is much less, about 0.05 liter per hour of swimming.<sup>28</sup> For chemicals with drinking water criteria, the proposed MCL of 1 ug/l was exceeded by beryllium in SW-4 (1.1 ug/l), and the Action Level of 15 ug/l was exceeded by lead in SW-2 (19 ug/l) and SW-4 (59.9 ug/l).<sup>29,30,31,32</sup> For these metals, carcinogenicity is a potential risk of concern because both are classified as Group B2 carcinogens. No EPA consensus for an oral CSF has been established for lead. For beryllium at this concentration, a 70-kilogram adult swimming seven days per year, two hours per day, for 30 years, ingesting 0.05 liter per hour, an increased oral cancer risk of approximately 6E-8 can be estimated.<sup>5</sup> Therefore, it can be stated that potential risks to swimmers in the river are due to the carcinogenicity of lead and beryllium; according to the no-threshold theory of carcinogenicity, any contact with carcinogens can increase overall cancer risk.

Sodium was detected in water from HW-1 at 21,900 ug/l, exceeding a guideline recommended by the American Heart Association (20,000 ug/l) to reduce the contribution of drinking water to total sodium intake.<sup>33</sup> While this can be important for persons on sodium-restricted diets, adverse health effects are not expected for the general population.

Report prepared by (b) (4)  
(b) (4), Toxicologist

Report reviewed by (b) (4)  
*[Signature]* (b) (4), DABT, Senior Toxicologist



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## **APPENDIX A**

Linfield Industrial Park

2598

**SUPPORT DOCUMENTATION FOR THE REVIEW OF  
ORGANIC ANALYSIS LAB DATA PACKAGE**

17794

Low organic

Company

3/90

(b) (4)

1992

APPLICABLE SAMPLE NO.: \_\_\_\_\_

Chy02-26

THE FOLLOWING TABLE INDICATES AREAS WHICH WERE EXAMINED IN DETAIL, THE IDENTIFIED PROBLEM AREAS, AND SUPPORT DOCUMENTATION ATTACHMENTS:

[illegible]

COMMENTS: \_\_\_\_\_

14744

## BLANK ANALYSIS RESULTS FOR TARGET COMPOUNDS

P.1

FRACTION	TYPE	CONC	MATRIX	SAMPLE #	SOURCE OF H <sub>2</sub> O	CONTAMINANTS (CONCENTRATION / DETECTION LIMIT)
VOA	low ag		lab blk	VBK26	lab	43 MeCl <sub>2</sub> 16.5 ug/l (1) 372 acetone 3.7 ug/l (2)
				VBK25		439 MeCl <sub>2</sub> 8.2 ug/l (1) 377 acetone 2.5 ug/l (2)
				VBK26		437 MeCl <sub>2</sub> 5.8 ug/l (1) 371 acetone 15.2 ug/l (2) 672 2-butanone 3.3 ug/l (2)
	low solid		lab blk	VBK27		421 MeCl <sub>2</sub> 27.5 ug/l 374 acetone 20.3 ug/l
				VBK28		423 MeCl <sub>2</sub> 39.0 ug/l (2) 377 acetone 18.6 ug/l (2)
				VBK29		442 MeCl <sub>2</sub> 13.5 ug/l (1) 383 acetone 19.2 ug/l (1)
				VBK34		422 MeCl <sub>2</sub> 25.6 ug/l (1) 376 acetone 27.1 ug/l (1)
	low ag		field blk	CH424	HMS	436 MeCl <sub>2</sub> 16.5 ug/l (1) 374 acetone 6.7 ug/l (2)
	low ag		trip blk	CH425	HMS	437 MeCl <sub>2</sub> 14.7 ug/l (1) 373 acetone 7.5 ug/l (2)
BOA	low ag		lab blk	SBK208	lab	1656 BBP/PhH 1.6 ug/l (2) 1750 DEHP 1.6 ug/l (2)
				SBK245		ND
				SBK295		ND
	low ag		field blk	CH424	HMS	1750 DEHP 1.2 ug/l (2)

LABORATORY REPORTED FIELD BLANK DATA IS COMPARED WITH THE SAMPLE DATA IN A TABULATION FORM WITHIN THE SAMPLE ANALYTICAL DATA SUMMARY. TENTATIVELY IDENTIFIED COMPOUNDS IN BLANKS ARE LISTED ON A SEPARATE FORM.

## COMMENTS:

(1) RESULT REPORTED BY LABORATORY AND CONFIRMED BY REVIEWER.

(2) RESULT INFERRED FROM QUANTITATION LIST, DIAGNOSTICS, CHROMATOGRAM AND/OR SPECTRA.

ND = none detected

17744

## BLANK ANALYSIS RESULTS FOR TARGET COMPOUNDS

B2

FRACTION	TYPE	CONC	MATRIX	SAMPLE #	SOURCE OF H <sub>2</sub> O	CONTAMINANTS (CONCENTRATION / DETECTION LIMIT)
BNA	low solid		lab blk	SBK96	lab	1514 DNEP 0.52 ug/l (2)
				SBK96		1386 DEP 0.25 ug/l (2) 1448 BAP 2.1 ug/l (2)
				SBK97		1200 DEP 0.25 ug/l (2) 1430 DNEP 0.50 ug/l (2) 1757 DEP 0.3 ug/l (2)
PST	low solid		lab blk	PB1408	lab	ND
	field		blk	chy24	HMS	ND
	SOLID		LAB BLK	PB162	lab	heptachlor 0.14 ug/kg (1) gamma-BHC 0.08 ug/kg endosulfan sulfate 0.52 ug/kg methoxychlor 0.36 ug/kg DDE 0.18 ug/kg (3) gamma-BHC 0.10 ug/kg (3) beta-BHC 0.41 ug/kg (3) endosulfan I 0.18 ug/kg (3) endosulfan II 0.32 ug/kg (3)
	low solid		lab blk	PB148		heptachlor 0.07 ug/kg (1) methoxychlor 0.61 ug/kg (1) endosulfan I 0.09 ug/kg (3) beta-BHC 0.35 ug/kg (3) endosulfan II 0.33 ug/kg (3)
	low solid		lab blk	PB173		heptachlor 0.08 ug/kg (1) Aniclor 1260 9.2 ug/kg (1) gamma-BHC 0.41 ug/kg (3)

DF=

9.0x33

LABORATORY REPORTED FIELD BLANK DATA IS COMPARED WITH THE SAMPLE DATA IN A TABULATION FORM WITHIN THE SAMPLE ANALYTICAL DATA SUMMARY. TENTATIVELY IDENTIFIED COMPOUNDS IN BLANKS ARE LISTED ON A SEPARATE FORM.

## COMMENTS:

(1) RESULT REPORTED BY LABORATORY AND CONFIRMED BY REVIEWER.

(2) RESULT INFERRED FROM QUANTITATION LIST, DIAGNOSTICS, CHROMATOGRAM AND/OR SPECTRA.

ND = none detected

\* = disagreement w/ amt. on 2 different columns.

(3) = compounds followed by (3) = peak only on column - most likely artifact ~~not used to question sample~~ NO corresponding peak on 2nd column.  
result is highest of two columns. (Not used to question sample) (3)



# BLANK ANALYSIS RESULTS FOR TENTATIVELY IDENTIFIED COMPOUNDS

ALL TENTATIVELY IDENTIFIED COMPOUNDS FOUND IN BLANK ANALYSES ARE LISTED BELOW:

1.0

SAMPLE #	FRACTION	SCAN # (S) OR R.T	ESTIMATED CONCENTRATION	COMPOUND NAME	COMMENTS
VBIKZB	VDA			ND	
VBIKAT	VDA			ND	
VBIKOL	VDA			ND	
VBIKTJ	VDA	488	2% ISO 618	MS	
"	"	1640	2% ISO 612	MS	(siloxane)
VBIKLV	VDA	441	3% ISO 610	MS	
	↓	1532	2% ISO 613	↓	(siloxane)
	↓	1643	" " "	↓	(siloxane)
VBIKLV	VDA			ND	
VBIKAY	VDA	334	1% ISO 610	MS	(furan)
"	"	490	6% ISO 610	MS	
CH424	VDA			ND	
CH425	VDA			ND	
SBKXOB	BNA	683	3 u/L	41, 45, 57, 87	
		689	8% ISO 772	MS	
		1656	5% ISO 1786	BB PTH.	
✓	✓	750	7% " "	✓	Delap
SBK95	BNA	451	2 u/L	main peak from 70-157	
		1082	3	siloxane	
		1197	2	43, 71	
		1500	6	43, 61, 69, 83, 97, 111	
		1509	5	43, 60, 101, 207	
		1975	3 ✓	69, 81, 95	
		547	7% ISO 641		
		930	5% ISO 837		
		1215	7% ISO 1122		
✓	✓	1333	4% ISO 1360		
SBK1295	BNA			ND	
CH424	BNA	684	5% ISO 772	MS	
	↓	690	5% " "	MS	
	↓	1750	5% ISO 1787	Delap	
SBK96	BNA	592	390 u/L	43, 69, 97, 112	FS: 653
		969	230	siloxane (73.54, 429...)	881
		1122	160	"	1169
		1258	160	"	1417
		1375	99	"	1865
		1660	66	"	2137
		1782	130	carbox acid (124)	
		2018	330 ✓	55, 57, 71, 83, 97, 112, 185	
		547	ISO 653	MS	
		609	5% ISO	1236 9% ISO 1169	MS
		612	5% ISO	1321 2% ISO 1417	
		672	3% ISO	1359 3% ISO	
		724	8% ISO	1400 1% ISO	
		825	4% ISO 881	1459 3% ISO	
		1206	3% ISO 1169	1460 3% ISO	
				1514 5% ISO	
				1574 4% ISO	
				1655 9% ISO 865	
				1740 5% ISO	
				1818 4% ISO	

ND = none detected  
 NLS = no library search conducted  
 ( ) = primary ans

PUB PTH.

17744

## BLANK ANALYSIS RESULTS FOR TENTATIVELY IDENTIFIED COMPOUNDS

ALL TENTATIVELY IDENTIFIED COMPOUNDS FOUND IN BLANK ANALYSES ARE LISTED BELOW:

p2

SAMPLE #	FRACTION	SCAN # (S) OR P.T	ESTIMATED CONCENTRATION	COMPOUND NAME	COMMENTS
Cont	BNA	1882	49% ISO 1865	NLS	
"	"	2029	8% ISO 2137	NLS	
SRLK86	BNA	653	20% U914	SHC	IS: 853
		756	1400	43, 69, 97	1049
		770	200	43, 69, 97, 112	1328
		717	5% ISO 853	NLS	1567
		729	6%		2041
		806	7%	THF?	2538
		835	5%		
		1001	1% ISO 1044		
		1112	" " "		
		1258	1% ISO 1528		
		1358	2% " "	PEP	
		1499	4% ISO 1567		
		1507	4%		
		1597	2%		
		1685	2%		
		1766	1%		
		1898	4% ISO 2041	BBP	
✓	✓	2263	4% " "		
✓	✓	827	4% ISO 853	✓	
SRLK91	BNA	565	230 U914	43, 69, 97	IS: 742
		657	270	43, 69, 97, 112	911
		703	130	43, 69, 97, 112	1154
		1064	170	1-bromine	1362
		620	5% ISO 142	NLS	1780
		634	5%		2208
		726	5%		
		788	3%		
		826	4%		
		907	1% ISO 911		
		971	3%	"	
		1097	8% ISO 1154		
		1200	2%	DEP	16
		1209	4%		
		1306	7% ISO 132		
		1376	3%		
		1389	6%		
		1430	4%	ONBP	17
		1464	5%		
		1544	4%		
		1548	2%		
		1611	6% ISO 1780		
		1635	4%		
		1726	4%		
1798		1759	9%	DEHP	1.0
		1881	3%		
		1979	2% ISO 2208		
		2102	6%	"	N/

ND = none detected

NLS = no library search conducted

( ) = primary ions

IS = internal standards

2A  
WATER VOLATILE SYSTEM MONITORING COMPOUND RECOVERY

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

	EPA SAMPLE NO.	SMC1 (TOL) #	SMC2 (BFB) #	SMC3 (DCE) #	OTHER	TOT OUT
01	CHY02	99	97	101	0	0
02	CHY03	101	97	103	0	0
03	CHY04	98	90	99	0	0
04	CHY05	97	89	101	0	0
05	CHY06	101	93	100	0	0
06	CHY07	100	86	103	0	0
07	CHY08	108	93	103	0	0
08	CHY09	102	109	98	0	0
09	CHY10	104	94	111	0	0
10	CHY24	106	93	110	0	0
11	CHY25	106	92	107	0	0
12	CHY07MS	99	90	114	0	0
13	CHY07MSD	102	89	103	0	0
14	VBLKZB	104	92	102	0	0
15	VBLKAJ	100	88	103	0	0
16	VBLKDC	106	110	99	0	0

QC LIMITS

SMC1 (TOL) = Toluene-d8 ( 88-110)  
 SMC2 (BFB) = Bromofluorobenzene ( 86-115)  
 SMC3 (DCE) = 1,2-Dichloroethane-d4 ( 76-114)

# Column to be used to flag recovery values  
 \* Values outside of contract required QC limits  
 D System Monitoring Compound diluted out

## WATER VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
 Matrix Spike - EPA Sample No.: CHY07

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)	MS CONCENTRATION (ug/L)	MS % REC #	QC LIMITS REC.
1,1-Dichloroethene	50.00	0	56.90	114	61-145
Trichloroethene	50.00	0	44.70	89	71-120
Benzene	50.00	0	47.60	95	76-127
Toluene	50.00	0	46.50	93	76-125
Chlorobenzene	50.00	0	44.80	90	75-130

COMPOUND	SPIKE ADDED (ug/L)	MSD CONCENTRATION (ug/L)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
1,1-Dichloroethene	50.00	63.40	127	11	14	61-145
Trichloroethene	50.00	43.70	87	2	14	71-120
Benzene	50.00	48.70	97	2	11	76-127
Toluene	50.00	45.10	90	3	13	76-125
Chlorobenzene	50.00	44.70	89	1	13	75-130

# Column to be used to flag recovery and RPD values with an asterisk

\* Values outside of QC limits

RPD: 0 out of 5 outside limits

Spike Recovery: 0 out of 10 outside limits

COMMENTS: CLP,17744,CHY02,CHY07,LOW,WATER,477460,VOA,EPA,F50051  
 DB624,CS920202A51,BF920202A51,CB920202A51,CN077447A51,CN077448A

2B  
SOIL VOLATILE SYSTEM MONITORING COMPOUND RECOVERY

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
 Level: (low/med) LOW

	EPA SAMPLE NO.	SMC1 (TOL) #	SMC2 (BFB) #	SMC3 (DCE) #	OTHER	TOT OUT
01	CHY11	110	103	107	0	0
02	CHY12	106	98	106	0	0
03	CHY13	108	94	110	0	0
04	CHY14	101	96	98	0	0
05	CHY15	112	101	110	0	0
06	CHY16	96	98	104	0	0
07	CHY17	127	111	118	0	0
08	CHY18	110	98	107	0	0
09	CHY19	105	95	103	0	0
10	CHY20	123	104	108	0	0
11	CHY21	119	99	104	0	0
12	CHY22	109	101	103	0	0
13	CHY23	93	90	96	0	0
14	CHY26	86	80	85	0	0
15	CHY18MS	109	98	105	0	0
16	CHY18MSD	108	97	107	0	0
17	VBLKT7	104	97	100	0	0
18	VBLKV8	93	89	89	0	0
19	VBLKW9	93	89	98	0	0
20	VBLKA4	101	96	98	0	0

QC LIMITS

SMC1 (TOL) = Toluene-d8 ( 84-138)  
 SMC2 (BFB) = Bromofluorobenzene ( 59-113)  
 SMC3 (DCE) = 1,2-Dichloroethane-d4 ( 70-121)

# Column to be used to flag recovery values  
 \* Values outside of contract required QC limits  
 D System Monitoring Compound diluted out

3B  
SOIL VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
 Matrix Spike - EPA Sample No.: CHY18 Level: (low/med) LOW

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC LIMITS REC.
1,1-Dichloroethene	69.40	0	68.19	98	59-172
Trichloroethene	69.40	0	63.33	91	62-137
Benzene	69.40	0	74.45	107	66-142
Toluene	69.40	0	69.03	99	59-139
Chlorobenzene	69.40	0	66.11	95	60-133

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
1,1-Dichloroethene	69.40	72.78	105	7	22	59-172
Trichloroethene	69.40	64.86	93	2	24	62-137
Benzene	69.40	76.11	110	3	21	66-142
Toluene	69.40	72.50	104	5	21	59-139
Chlorobenzene	69.40	69.17	100	5	21	60-133

# Column to be used to flag recovery and RPD values with an asterisk

\* Values outside of QC limits

RPD: 0 out of 5 outside limits  
 Spike Recovery: 0 out of 10 outside limits

COMMENTS: CLP  
 CAP, GT920131C18, BG920131C18, , , ,

2D  
SOIL SEMIVOLATILE SURROGATE RECOVERY

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
 Level: (low/med) LOW

	EPA SAMPLE NO.	S1 (NBZ) #	S2 (FBP) #	S3 (TPH) #	S4 (PHL) #	S5 (2FP) #	S6 (TBP) #	S7 (2CP) #	S8 (DCB) #	TOT OUT
01	CHY11	88	83	67	70	72	45	57	78	0
02	CHY12	58	51	76	47	42	33	35	48	0
03	CHY13	62	63	70	50	50	36	40	60	0
04	CHY14	75	64	79	66	69	39	54	60	0
05	CHY15	0 D	0 D	87	0 D	0 D	64	0 D	0 D	0
06	CHY16	58	58	76	38	19 *	26	24	34	1
07	CHY17	76	68	79	60	47	13 *	40	63	1
08	CHY18	79	64 ✓	79 ✓	59 ✓	48 ✓	18 *	42	63	1
09	CHY18DL	84	75 ✓	91 ✓	69 ✓	68 ✓	46 ✓	55	68	0
10	CHY19	77	71	66	64	51	21	43	64	0
11	CHY20	45	59	63	45	66	27	45	55	0
12	CHY20DL	71	65	87	45	56	46	41	0 D	0
13	CHY21	42	44	40	42	46	40	34	41	0
14	CHY22	80	71	72	50	28	2 *	27	64	1
15	CHY22RE	58	71	81	55	35 ✓	8 *	32 ✓	53	1
16	CHY23	47	45	49	37	39	34	31	43	0
17	CHY26	72	65	70	68	62	55	50	62	0
18	CHY18MS	71	60	72	51	40	12 *	35	60	1
19	CHY18MSD	79	69	82	66	63	53 ✓	49	69	0
20	SBLK96	61	52	88	47	44	37	35	52	0
21	SBLK86	38	39	44	41	41	23	30	40	0
22	SBLK97	51	49	58	48	45	42	37	47	0

QC LIMITS

S1 (NBZ) = Nitrobenzene-d5 ( 23-120)  
 S2 (FBP) = 2-Fluorobiphenyl ( 30-115)  
 S3 (TPH) = Terphenyl-d14 ( 18-137)  
 S4 (PHL) = Phenol-d5 ( 24-113)  
 S5 (2FP) = 2-Fluorophenol ( 25-121)  
 S6 (TBP) = 2,4,6-Tribromophenol ( 19-122)  
 S7 (2CP) = 2-Chlorophenol-d4 ( 20-130) (advisory)  
 S8 (DCB) = 1,2-Dichlorobenzene-d4 ( 20-130) (advisory)

# Column to be used to flag recovery values  
 \* Values outside of contract required QC limits  
 D Surrogate diluted out

&amp;l0o61f6D

3D

## SOIL SEMIVOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPUCase No.: 17744SAS No.: 6579HQSDG No.: CHY11Matrix Spike - EPA Sample No.: CHY18Level: (low/med) LOW

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC LIMITS REC.
Phenol	3430	0	1928	56	26- 90
2-Chlorophenol	3430	0	1704	50	25-102
1,4-Dichlorobenzene	2280	0	1366	60	28-104
N-Nitroso-di-n-prop. (1)	2280	0	1467	64	41-126
1,2,4-Trichlorobenzene	2280	0	1435	63	38-107
4-Chloro-3-methylphenol	3430	0	1869	54	26-103
Acenaphthene	2280	88.65	1499	62	31-137
4-Nitrophenol	3430	0	2367	69	11-114
2,4-Dinitrotoluene	2280	0	1307	57	28- 89
Pentachlorophenol	3430	0	78.14	2 *	17-109
Pyrene	2280	2691	3263	25 *	35-142

TBP surr. low

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
Phenol	3430	2276	66	16	35	26- 90
2-Chlorophenol	3430	2166	63	23	50	25-102
1,4-Dichlorobenzene	2280	1426	63	5	27	28-104
N-Nitroso-di-n-prop. (1)	2280	1604	70	9	38	41-126
1,2,4-Trichlorobenzene	2280	1471	65	3	23	38-107
4-Chloro-3-methylphenol	3430	2490	73	30	33	26-103
Acenaphthene	2280	1650	68	9	19	31-137
4-Nitrophenol	3430	2495	73	6	50	11-114
2,4-Dinitrotoluene	2280	1353	59	3	47	28- 89
Pentachlorophenol	3430	302.0	9 *	127 *	47	17-109
Pyrene	2280	3564	38 ✓	41 *	36	35-142

TBP surr. ch

(1) N-Nitroso-di-n-propylamine

low pyrene recov. possible due to sp1  
inhomogeneity -  
res. est.

# Column to be used to flag recovery and RPD values with an asterisk

\* Values outside of QC limits

RPD: 2 out of 11 outside limitsSpike Recovery: 3 out of 22 outside limits

COMMENTS: CLP

CAP, HH920210C21, DH920210C21, , , ,



2C  
WATER SEMIVOLATILE SURROGATE RECOVERY

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

	EPA SAMPLE NO.	S1 (NBZ) #	S2 (FBP) #	S3 (TPH) #	S4 (PHL) #	S5 (2FP) #	S6 (TBP) #	S7 (2CP) #	S8 (DCB) #	TOT OUT
01	CHY02	102	98	103	96	98	92	75	92	0
02	CHY03	87	81	103	79	83	77	61	81	0
03	CHY04	79	72	92	70	73	63	56	76	0
04	CHY05	90	82	109	79	78	75	63	86	0
05	CHY06	71	63	79	60	57	63	47	64	0
06	CHY07	83	71	94	75	79	74	60	78	0
07	CHY08	83	71	85	72	77	79	57	75	0
08	CHY09	84	73	70	72	79	72	57	84	0
09	CHY10	83	73	89	70	70	64	56	77	0
10	CHY24	65	65	93	73	70	53	55	64	0
11	CHY07MS	83	75	104	78	77	84	56	71	0
12	CHY07MSD	82	73	93	70	75	79	54	71	0
13	SBLK95	76	68	94	71	73	31	55	71	0
14	SBLK08	56	57	77	64	62	45	49	55	0
15	SBLK95	102	93	123	103	104	72	77	92	0

QC LIMITS

S1 (NBZ) = Nitrobenzene-d5 ( 35-114)  
S2 (FBP) = 2-Fluorobiphenyl ( 43-116)  
S3 (TPH) = Terphenyl-d14 ( 33-141)  
S4 (PHL) = Phenol-d5 ( 10-110)  
S5 (2FP) = 2-Fluorophenol ( 21-110)  
S6 (TBP) = 2,4,6-Tribromophenol ( 10-123)  
S7 (2CP) = 2-Chlorophenol-d4 ( 33-110) (advisory)  
S8 (DCB) = 1,2-Dichlorobenzene-d4 ( 16-110) (advisory)

# Column to be used to flag recovery values  
\* Values outside of contract required QC limits  
D Surrogate diluted out

3C  
WATER SEMIVOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
 Matrix Spike - EPA Sample No.: CHY07

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)	MS CONCENTRATION (ug/L)	MS % REC #	QC LIMITS REC.
Phenol	75.00	0	66.20	88	12-110
2-Chlorophenol	75.00	0	64.50	86	27-123
1,4-Dichlorobenzene	50.00	0	38.30	77	36- 97
N-Nitroso-di-n-prop. (1)	50.00	0	46.50	93	41-116
1,2,4-Trichlorobenzene	50.00	0	41.00	82	39- 98
4-Chloro-3-methylphenol	75.00	0	71.40	95	23- 97
Acenaphthene	50.00	0	43.00	86	46-118
4-Nitrophenol	75.00	1.320	92.10	121 *	10- 80
2,4-Dinitrotoluene	50.00	0	53.00	106 *	24- 96
Pentachlorophenol	75.00	0	39.00	52	9-103
Pyrene	50.00	0	54.30	109	26-127

COMPOUND	SPIKE ADDED (ug/L)	MSD CONCENTRATION (ug/L)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
Phenol	75.00	62.70	84	5	42	12-110
2-Chlorophenol	75.00	62.40	83	4	40	27-123
1,4-Dichlorobenzene	50.00	35.60	71	8	28	36- 97
N-Nitroso-di-n-prop. (1)	50.00	46.60	93	0	38	41-116
1,2,4-Trichlorobenzene	50.00	39.60	79	4	28	39- 98
4-Chloro-3-methylphenol	75.00	65.90	88	8	42	23- 97
Acenaphthene	50.00	40.50	81	6	31	46-118
4-Nitrophenol	75.00	87.10	114 *	6	50	10- 80
2,4-Dinitrotoluene	50.00	48.70	97 *	9	38	24- 96
Pentachlorophenol	75.00	64.70	86	49	50	9-103
Pyrene	50.00	48.20	96	13	31	26-127

(1) N-Nitroso-di-n-propylamine

# Column to be used to flag recovery and RPD values with an asterisk  
 \* Values outside of QC limits

RPD: 0 out of 11 outside limits  
 Spike Recovery: 4 out of 22 outside limits

COMMENTS: CLP  
 CAP, HG920203B52, DH920203B52, , ,

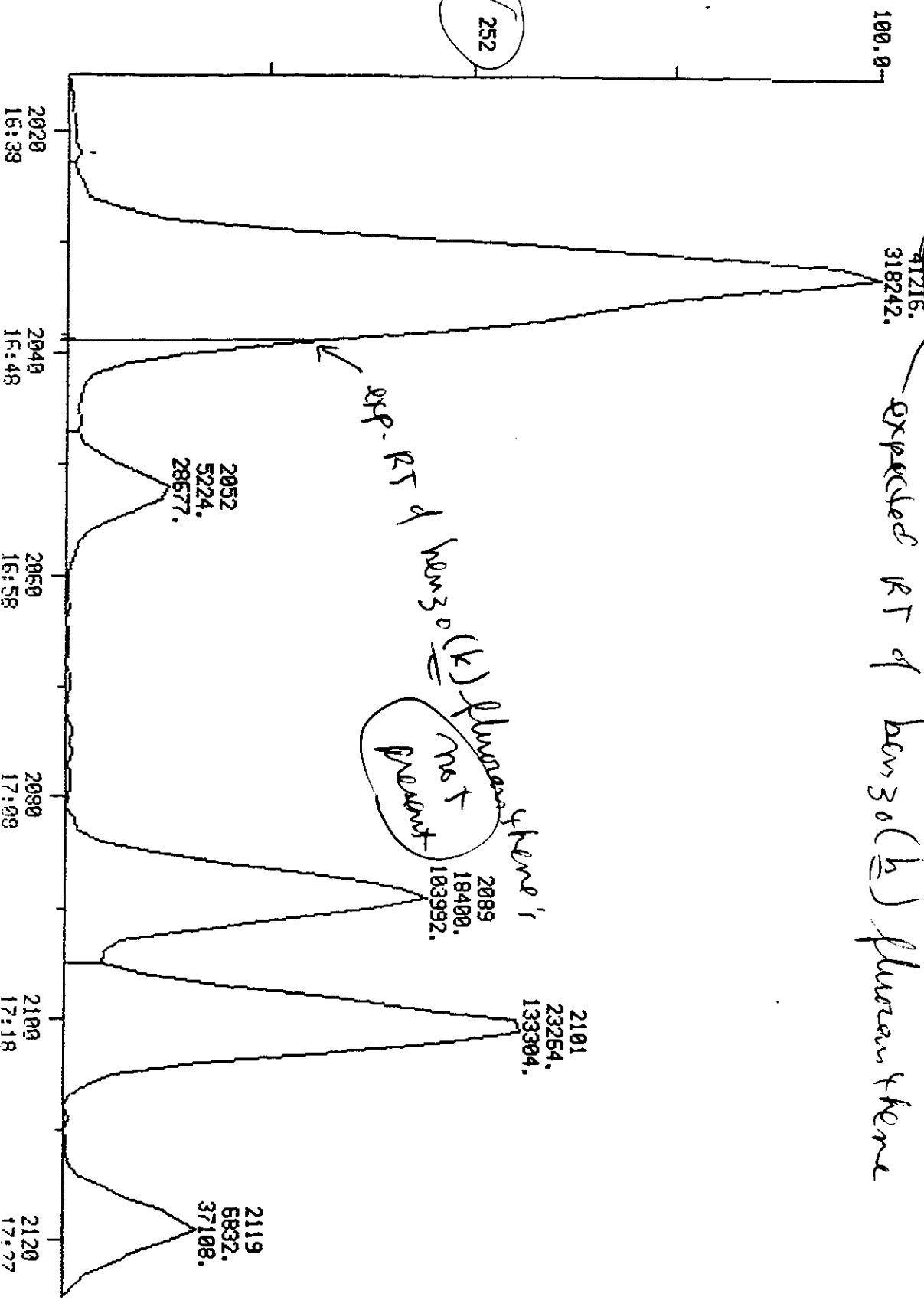
FORM III SV-1

Sample Data Package

17744 CHY02

3/90 CEE  
 2/21/96  
 251 A

MIDMASS CHROMATOGRAM  
 02/12/92 21:32:00  
 SAMPLE: 2UL CC#477472 ID#CHY11 RE  
 COND5.: EXTRACTED 2/4/92 1: 2 DILUTION  
 RANGE: G 1.2835 LABEL: H 1, 3.0 QUAN: A 1, 0.3 J 0 BASE: U 20, 3  
 DATA: GDUJ77472A21 #1  
 CALL: GDUJ77472A21 #2  
 CS#17744-CHY11  
 OMR21  
 SCANS 2015 TO 2125



SAMPLE DATA PACKAGE  
 252.075  
 ± 0.500

17744 CHY 11

475

51 444 PHENANTHRENE (Q4#7) <85-01-8>  
 52 583 CARBAZOLE  
 53 403 ANTHRACENE (Q4#8) <120-12-7>  
 54 426 DI-N-BUTYL PHTHALATE (Q4#9) <84-74-2>  
 55 431 FLUORANTHENE (Q4#10) <206-44-0>  
 56 \*459 D12-CHRYSENE (IS#5)  
 57 445 PYRENE (Q5#3) <129-00-0>  
 58 415 BUTYLBENZYL PHTHALATE (Q5#4) <85-68-7>  
 59 423 3,3'-DICHLOROBENZIDINE (Q5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (Q5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (Q5#7) <117-81-7>  
 62 418 CHRYSENE (Q5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (Q6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (Q6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (Q6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (Q6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (Q6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (Q6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (Q6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CHY 11

Review added  
 Benz (c)  
 used popn RPF

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1403	11:33	45	1.003	A BV	318796.	26.529 NG	4.33 y
52	167	1440	11:52	45	1.029	A BB	22824.	2.666 NG	0.44 y
53	178	1411	11:37	45	1.009	A VV	62644.	5.992 NG	0.98 y
54	NOT FOUND								
55	202	1403	13:12	45	1.146	A BV	532992.	60.866 NG	9.93 y
56	240	1846	15:12	56	1.000	A VB	157164.	20.000 NG	3.26 y
57	202	1641	13:31	56	0.889	A BB	477385.	30.897 NG	5.04 y
58	149	1751	14:25	56	0.949	A BB	11723.	1.428 NG	0.23 y
59	NOT FOUND								
	228	1845	15:12	56	0.999	A BV	229611.	25.381 NG	4.14 y
60	149	1850	15:14	56	1.002	A BB	37186.	2.979 NG	0.49 y
62	228	1851	15:15	56	1.003	A VV	173006.	19.377 NG	3.16 y
63	264	2113	17:24	63	1.000	A BB	129628.	20.000 NG	3.26 y
64	NOT FOUND								
65	252	2033	16:45	63	0.962	A BV	322113.	39.375 NG	7.05 y

x 92 = 4,000 ug/kg  
 (34%)  
 yes  
 40.344  
 20.047  
 13.077  
 12.505  
 4.439  
 6(b)f

Quantitation Report File: GDJ77472A21

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
55	252	2033	16:45	63	0.962	A BV	322113.	39.375 NG	6.42 y
57	252	2101	17:18	63	0.994	A VV	144998.	21.763 NG	3.55 y
68	276	2411	19:51	63	1.141	A BB	105072.	14.414 NG	2.35 y
69	NOT FOUND								
70	276	2504	20:37	63	1.185	A BB	86434.	13.337 NG	2.18 y
71	112	486	4:00	1	0.728	A BB	157976.	27.140 NG	4.43 y
72	99	618	5:05	1	0.925	A BV	192244.	26.368 NG	4.30 y
73	132	636	5:14	1	0.952	A BB	145932.	21.386 NG	3.49 y
74	152	691	5:41	1	1.034	A BB	88820.	19.615 NG	3.20 y
75	82	753	6:12	13	0.871	A BV	161808.	22.017 NG	3.59 y
76	172	1041	8:34	26	0.904	A BV	231208.	20.846 NG	3.40 y
77	330	1282	10:33	26	1.114	A BB	200000.	20.000 NG	3.26 y

02/13/12

55 431 FLUORANTHENE (G4#10) <206-44-0>  
 56 \*459 D12-CHRYSENE (IS#5)  
 57 445 PYRENE (G5#3) <129-00-0>  
 58 415 BUTYLBENZYL PHTHALATE (G5#4) <85-68-7>  
 59 423 3,3'-DICHLOROBENZIDINE (G5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (G5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (G5#7) <117-81-7>  
 62 418 CHRYSENE (G5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (G6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (G6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (G6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (G6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (G6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (G6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (G6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CC  
 STD  
 2/12

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1407	11:35	45	1.003	A BV	402592.	25.000 NG	1.30
52	167	1443	11:53	45	1.029	A BV	286820.	25.000 NG	1.30
53	178	1415	11:39	45	1.009	A VB	350268.	25.000 NG	1.30
54	149	1502	12:22	45	1.071	A BV	551675.	25.000 NG	1.30
55	202	1606	13:14	45	1.145	A BV	293372.	25.000 NG	1.30
56	240	1849	15:14	56	1.000	A BB	121620.	20.000 NG	1.04
57	202	1644	13:32	56	0.889	A BB	298912.	25.000 NG	1.30
58	149	1754	14:27	56	0.949	A BV	158868.	25.000 NG	1.30
59	252	1843	15:11	56	0.997	A BB	23160.	25.000 NG	1.30
60	228	1848	15:13	56	0.999	A BV	175017.	25.000 NG	1.30
61	149	1853	15:16	56	1.002	A BB	241512.	25.000 NG	1.30
62	228	1854	15:16	56	1.003	A VB	172728.	25.000 NG	1.30
63	264	2114	17:25	63	1.000	A BB	83140.	20.000 NG	1.04
64	149	1963	16:10	63	0.929	A BV	318104.	25.000 NG	1.30
65	252	2034	16:45	63	0.962	A BV	119496.	25.000 NG	1.30

is 5  
 is 6  
 b(k)f

SAMPLE DATA PACKAGE

17744 CHY 11

1759

Quantitation Report File: HG920212B21

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2039	16:47	63	0.965	A VB	131170.	25.000 NG	1.30
67	252	2102	17:19	63	0.994	A BV	106828.	25.000 NG	1.30
68	276	2411	19:51	63	1.140	A BB	116884.	25.000 NG	1.30
69	278	2419	19:55	63	1.144	A BB	95000.	25.000 NG	1.30
70	276	2500	20:35	63	1.183	A BB	103912.	25.000 NG	1.30
71	112	487	4:01	1	0.725	A BB	157576.	25.000 NG	1.30
72	99	621	5:07	1	0.924	A BV	197372.	25.000 NG	1.30
73	132	639	5:16	1	0.951	A BV	184728.	25.000 NG	1.30
74	152	695	5:43	1	1.034	A BB	122584.	25.000 NG	1.30
75	82	757	6:14	13	0.871	A BB	228712.	25.000 NG	1.30
76	172	1045	8:36	26	0.905	A BB	326200.	25.000 NG	1.30
77	330	1286	10:35	26	1.113	A BB	56208.	25.000 NG	1.30
78	244	1671	13:46	56	0.904	A BB	196454.	25.000 NG	1.30

b(k)f

No Ret(L) Ratio RRT(L) Ratio Amt+ Amt+ (L) ...

COMPUCHEM LABS, INC.

MID LIBRARY SEARCH (LIBRARYINC)

DATA: GDU77472A21 #2504

ENHANCED (128 2N 0T)

BASE M/Z: 67  
R/C: 12254.

02/12/92 21:32:22 + 20137  
SAMPLE: 2UL OS#477472 ID#CHY11 RE  
CONVS: 1 EXTRACTED 2/4/92 11 2 COLLECTION

OS#17744-CHY11 CHM21

SAMPLE

5224

C22, H12

5094

M WT 750  
R PK 275

RANK 1

# 79

PUR 145

408 BENZO(C,H,1)PERYLENE (06#8) (191-24-2)

SAMPLE MINUS LIBRARY

5094

M/Z

50

100

150

200

250

Strong as  
tentative

not characteristic  
mass profile

SAMPLE DATA PACKAGE

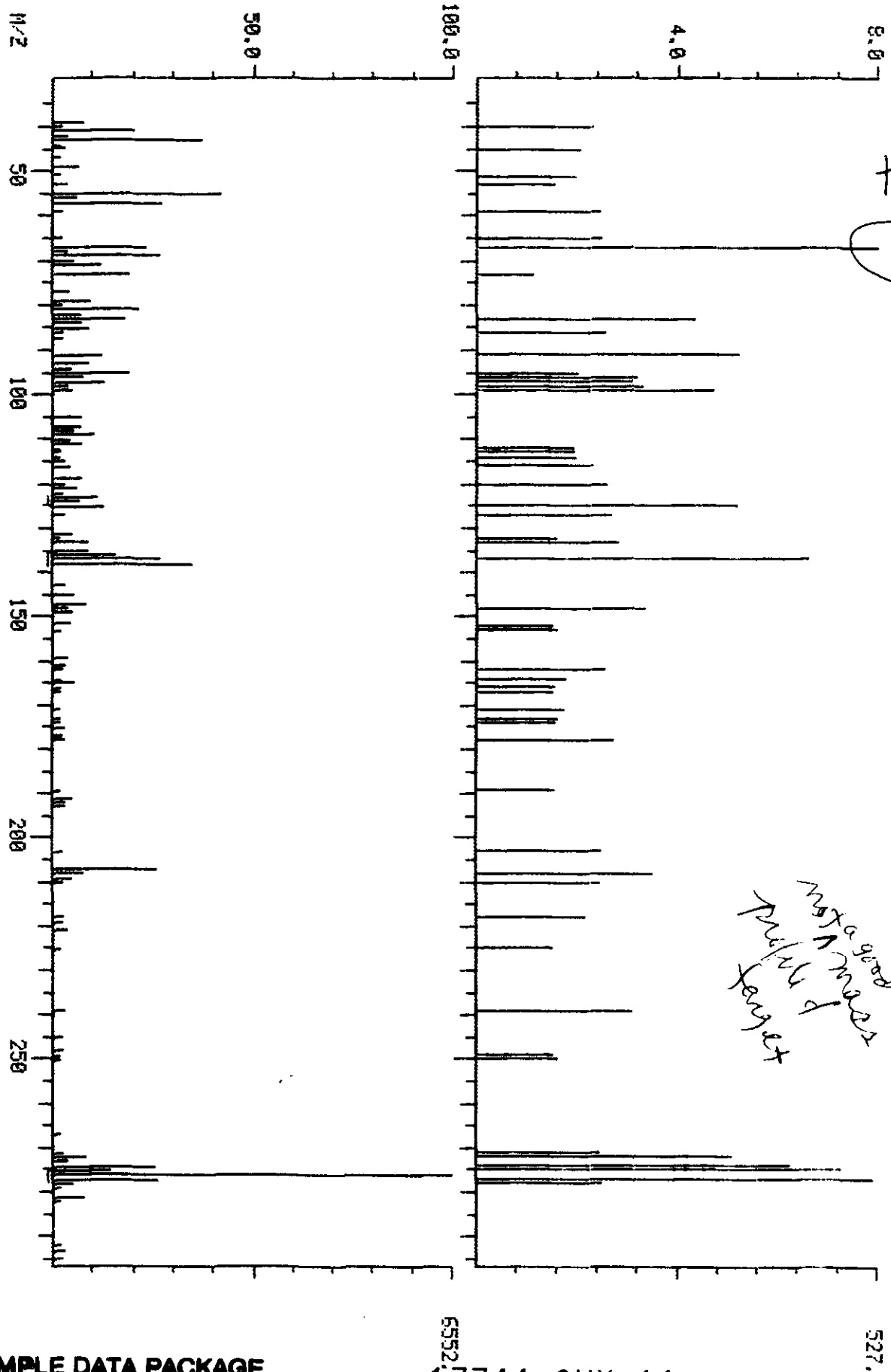
17744 CHY 11

482

MID DUAL MASS SPECTRUM  
02/12/92 21:32:00 + 20:37  
SECOND SPECTRUM  
SAMPLE: 2UL CC#477472 ID#CHY11 BE  
408 BENZO(G,H,I)PERYLENE (06#8) (191-24-2)

DATA: GDU77472021 #2504 BASE M/Z: 67/ 276  
ENHANCED (108 2N 0T) RIC: 10159. / 63359.  
DATA: UNENHANCED #2504 COMPUCHEN LABS, INC.  
01A021

*not a good  
match  
with  
target*

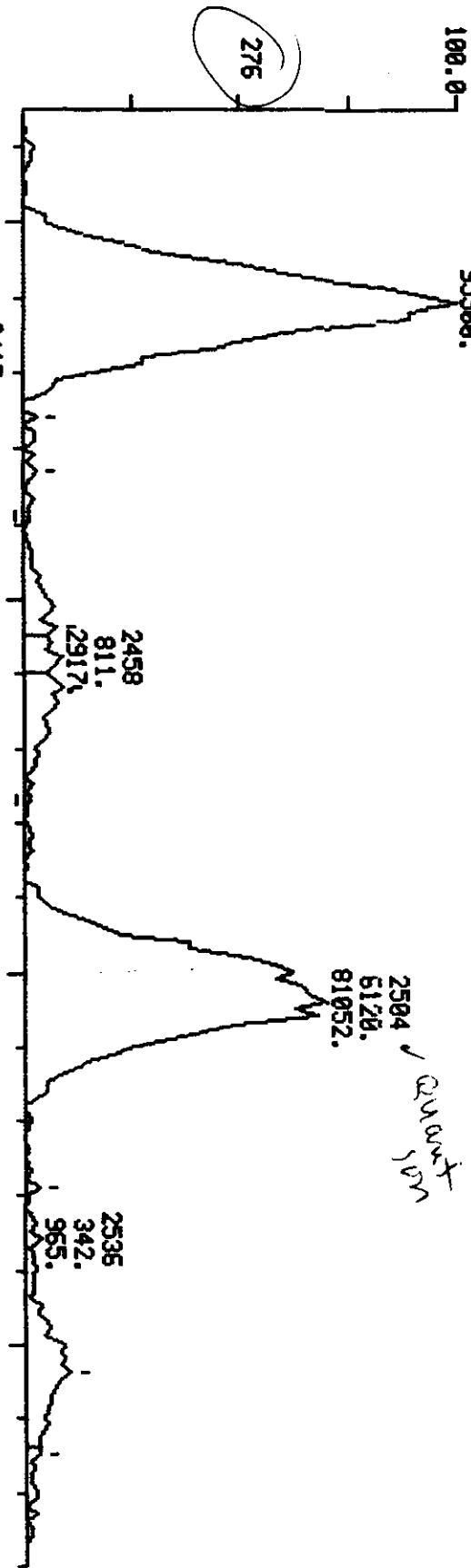


MIDMASS CHROMATOGRAMS  
 02/12/92 21:32:00  
 SAMPLE: 20L CC#477472 ID#CHY11 RE  
 COND.: EXTRACTED 2/4/92 1: 2 DILUTION  
 RANGE: G 1, 2855 LABEL: N 1, 3.0 QUAN: A 1, 0.3 J 0 BASE: U 20, 3

DATA: GDJ77472A21 #1

CALL: GDJ77472A21 #2  
 CS#17744-CHY11

SCANS 2385 TO 2580  
 OMR21

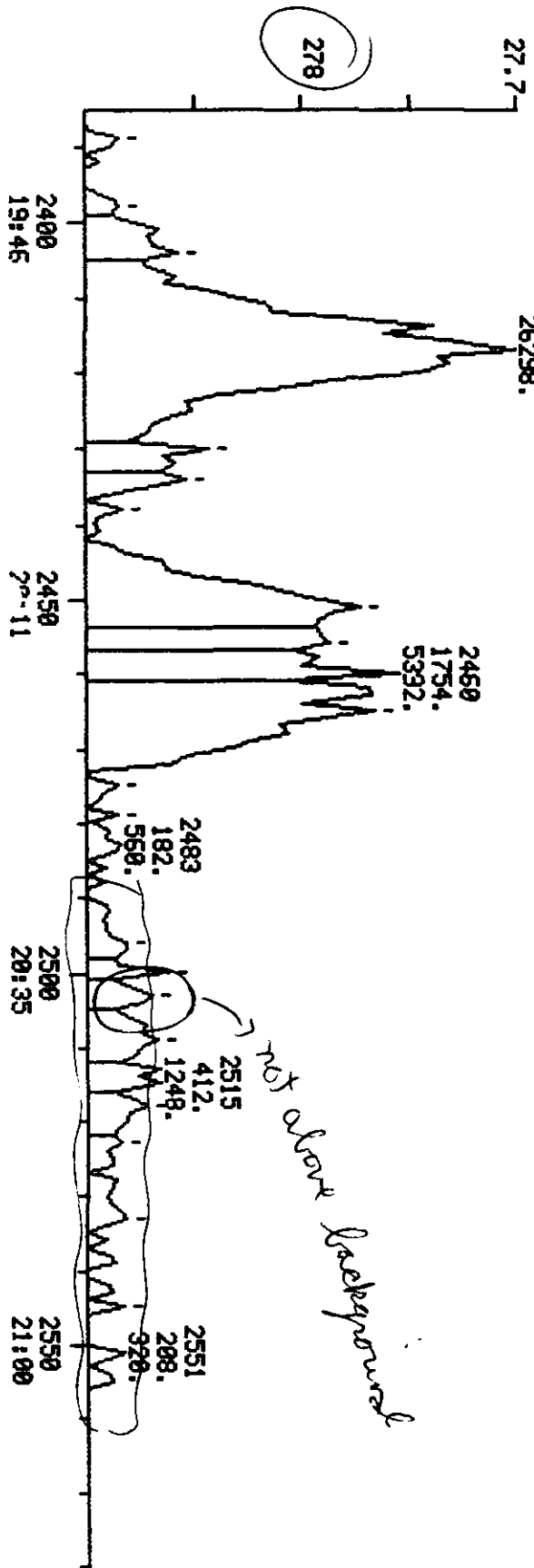


276.083  
 ± 0.500

8720.

17744 CHY 11

484



278.083  
 ± 0.500

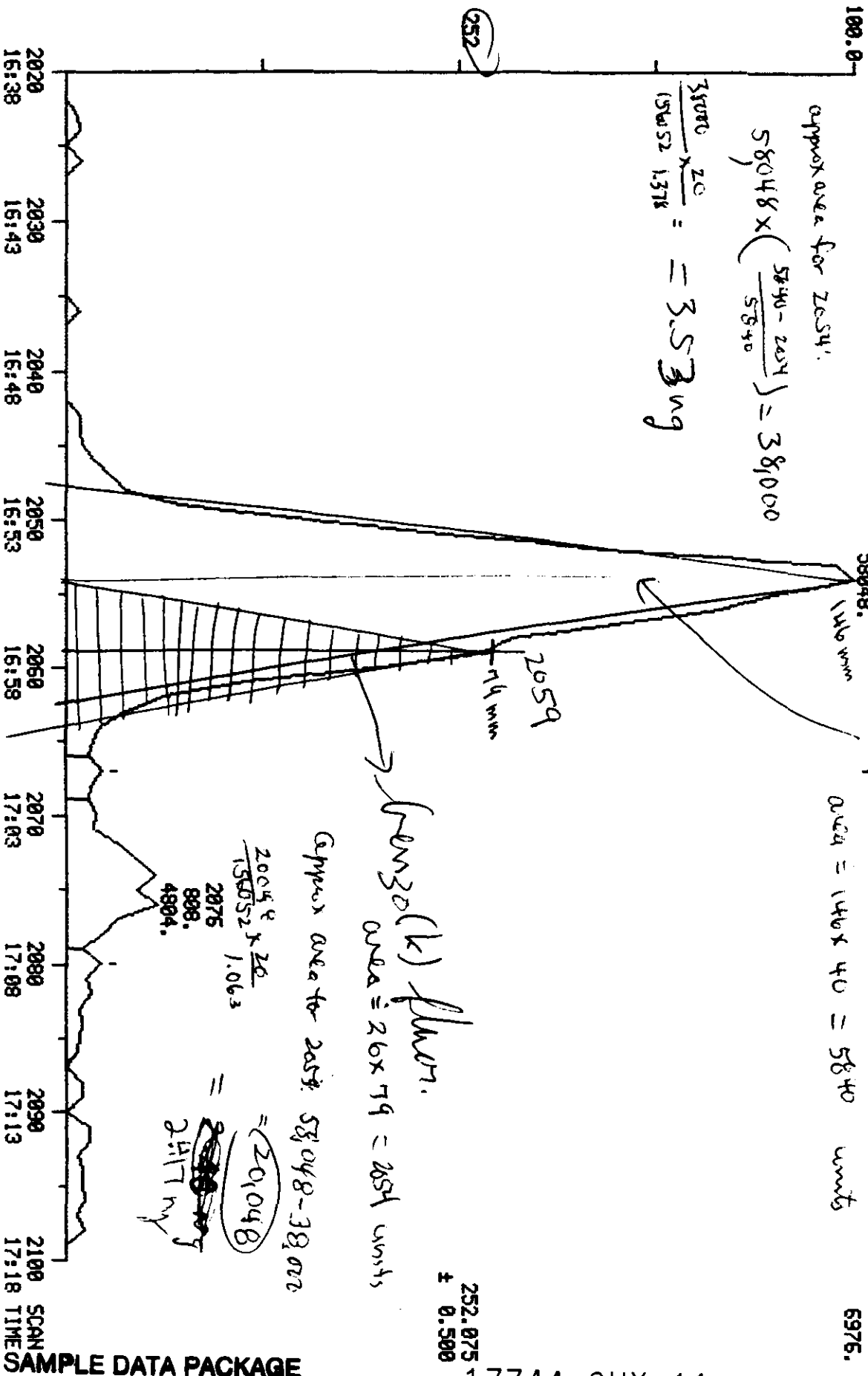
SCAN  
 TIME

SAMPLE DATA PACKAGE



16.12-52

MIDMASS CHROMATOGRAM  
 02/10/92 15:30:00  
 SAMPLE: 2UL CC#477476 ID#CHY12 RE  
 COND: : EXTRACTED 2/4/92 UNOILUTED  
 RANGE: C 1,2855 LABEL: N 1, 4.0 QUANT: 2, 1.0 J 0 BASE: U 20, 3  
 DATA: GR077476C21 #1 SCANS 2020 TO 2100  
 CALL: GR077476C21 #2  
 CS#17744-CHY11 OMR21



SAMPLE DATA PACKAGE

17744 CHY 11

52 583 CARBAZOLE  
 53 403 ANTHRACENE (Q4#8) <120-12-7>  
 54 426 DI-N-BUTYL PHTHALATE (Q4#9) <84-74-2>  
 55 431 FLUORANTHENE (Q4#10) <206-44-0>  
 56 \*459 D12-CHRYSENE (IS#5)  
 57 445 PYRENE (Q5#3) <129-00-0>  
 58 415 BUTYLBENZYL PHTHALATE (Q5#4) <85-68-7>  
 59 423 3,3'-DICHLOROBENZIDINE (Q5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (Q5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (Q5#7) <117-81-7>  
 62 418 CHRYSENE (Q5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (Q6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (Q6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (Q6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (Q6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (Q6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (Q6#7) <53-70-3>  
 70 408 BENZO(Q,H,I)PERYLENE (Q6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

Ch412

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1420	11:42	45	1.003	A BV	38952.	2.013 NG	0.50 yes
52	NOT FOUND								
53	NOT FOUND								
54	149	1515	12:29	45	1.070	A BB	21156.	1.079 NG	0.27 yes
55	202	1620	13:20	45	1.144	A BV	75756.	4.987 NG	1.23 yes
56	240	1864	15:21	56	1.000	A BB	247296.	20.000 NG	4.92 155
57	202	1659	13:40	56	0.890	A BB	71364.	4.397 NG	1.08 yes
58	149	1768	14:34	56	0.948	A BB	6620.	1.037 NG	0.26 yes
59	NOT FOUND								
60	228	1858	15:23	56	1.002	A*BV	3620 71176.	2.912 5.724 NG	1.41 yes
61	149	1866	15:22	56	1.001	A BB	49501.	4.829 NG	1.19 yes
62	228	1868	15:23	56	1.002	A VV	33032 84832.	2.866 2.978 NG	0.73 yes
63	264	2138	17:36	63	1.000	A BV	156052.	20.000 NG	4.92 156
64	NOT FOUND								
65	252	2054	16:55	63	0.961	A BV	381000 57478.	3.534 = 260 5.347 NG	1.32 yes

# SAMPLE DATA PACKAGE

17744 CHY 11

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2054	16:55	63	0.961	A BV	37045 57478.	6.927 NG	1.70 yes
67	252	2125	17:30	63	0.994	A VB	20160.	2.499 NG	0.61 yes
68	276	2443	20:07	63	1.143	A*BB	11004.	1.214 NG	0.30 yes
69	NOT FOUND								
70	276	2540	20:55	63	1.188	A BB	9116.	1.190 NG	0.29 yes
71	112	501	4:08	1	0.734	A BV	221964.	31.843 NG	7.83
72	99	632	5:12	1	0.925	A BV	291655.	35.430 NG	8.72
73	132	651	5:22	1	0.953	A BB	212764.	26.408 NG	6.50
74	152	706	5:49	1	1.034	A BB	134212.	24.090 NG	5.93
75	82	768	6:19	13	0.873	A BV	233100.	29.205 NG	7.19
76	172	1057	8:42	26	0.905	A BB	428016.	25.675 NG	6.32
77	330	1299	10:42	26	1.112	A BB	72364.	24.454 NG	6.02
78	244	1685	13:53	56	0.904	A BB	423108.	38.075 NG	9.37

No Ret(L) Ratio RRT(L) Ratio Amnt Amnt(L) R. Fac R Fac(L) Ratio

52 383 CARBAZOLE  
 53 403 ANTHRACENE (G4#8) <120-12-7>  
 54 426 DI-N-BUTYL PHTHALATE (G4#9) <84-74-2>  
 55 431 FLUORANTHENE (G4#10) <206-44-0>  
 56 \*459 D12-CHRYSENE (IS#5)  
 57 445 PYRENE (G5#3) <129-00-0>  
 58 415 BUTYLBENZYL PHTHALATE (G5#4) <85-68-7>  
 59 423 3,3'-DICHLOROBENZIDINE (G5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (G5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (G5#7) <117-81-7>  
 62 418 CHRYSENE (G5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (G6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (G6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (G6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (G6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (G6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (G6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (G6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

STD  
 2/10

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1421	11:42	45	1.003	A BV	342912.	25.000 NG	1.30
52	167	1459	12:01	45	1.030	A BV	167213.	25.000 NG	1.30
53	178	1430	11:47	45	1.009	A VB	266276.	25.000 NG	1.30
54	149	1516	12:29	45	1.070	A BB	599981.	25.000 NG	1.30
55	202	1622	13:21	45	1.145	A BB	291538.	25.000 NG	1.30
56	240	1865	15:21	56	1.000	A BB	200340.	20.000 NG	1.04
57	202	1660	13:40	56	0.890	A BB	353288.	25.000 NG	1.30
58	149	1769	14:34	56	0.949	A BV	127104.	25.000 NG	1.30
59	252	1859	15:19	56	0.997	A BB	24716.	25.000 NG	1.30
60	228	1864	15:21	56	0.999	A BV	275795.	25.000 NG	1.30
61	149	1867	15:22	56	1.001	A BV	291028.	25.000 NG	1.30
62	228	1870	15:24	56	1.003	A VB	208441.	25.000 NG	1.30
63	264	2138	17:36	63	1.000	A BV	113516.	20.000 NG	1.04
64	149	1978	16:17	63	0.925	A VB	249340.	25.000 NG	1.30
65	252	2055	16:55	63	0.961	A BV	203783.	25.000 NG	1.30

# SAMPLE DATA PACKAGE

17744 CHY 11

1739

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2060	16:58	63	0.964	A VB	167083.	25.000 NG	1.30
67	252	2126	17:30	63	0.994	A BB	133588.	25.000 NG	1.30
68	276	2448	20:10	63	1.145	A BB	133551.	25.000 NG	1.30
69	278	2455	20:13	63	1.148	A BB	123552.	25.000 NG	1.30
70	276	2544	20:57	63	1.190	A BB	128353.	25.000 NG	1.30
71	112	499	4:07	1	0.730	A BB	152868.	25.000 NG	1.30
72	99	632	5:12	1	0.924	A BV	182372.	25.000 NG	1.30
73	132	651	5:22	1	0.952	A BB	148420.	25.000 NG	1.30
74	152	707	5:49	1	1.034	A BB	108440.	25.000 NG	1.30
75	82	770	6:20	13	0.873	A VB	207772.	25.000 NG	1.30
76	172	1058	8:43	26	0.905	A BB	250424.	25.000 NG	1.30
77	330	1300	10:42	26	1.112	A BB	59392.	25.000 NG	1.30
78	244	1685	13:53	56	0.903	A BB	215176.	25.000 NG	1.30

No Ret(L) Ratio RRT(L) Ratio Amt Amt/L

MIDMASS CHROMATOGRAM  
02/12/92 22:10:00  
SAMPLE: 2UL CC#477483 10#CHY13 RE  
CONDS.: EXTRACTED 2/4/92 UNOILUTED  
RANGE: G 1.2855 LABEL: N 1.3.0 QUAN: A 1.0.3 J 0 BASE: U 20, 3

DATA: G2J77483R21 #1 SCANS 2015 TO 2125  
CALL: G2J77483R21 #2  
CS#17744-CHY11 OMA21

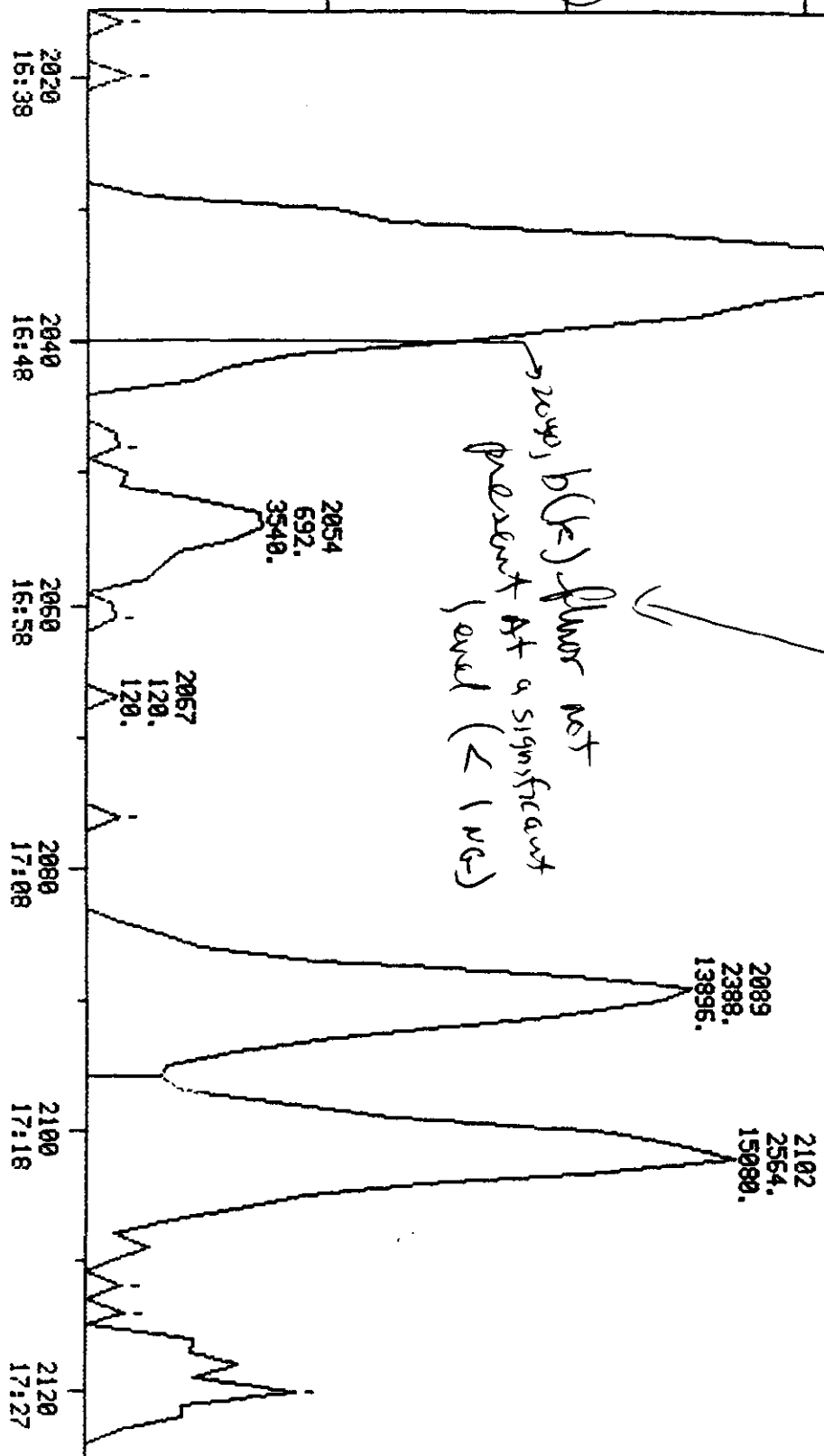
100.0

3772.  
28332.

*benz (b) fluor.*

*pred. b(k) = 2040*

*2040, b(k) fluor not present at a significant level (< 1 ng)*



252.075  
± 0.500

3772.

55 431 FLUORANTHENE (G4#10) <206-44-0>  
 56 \*459 D12-CHRYSENE (IS#5)  
 57 445 PYRENE (G5#3) <129-00-0>  
 58 415 BUTYLBENZYL PHTHALATE (G5#4) <85-68-7>  
 59 423 3,3'-DICHLOROBENZIDINE (G5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (G5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (G5#7) <117-81-7>  
 62 418 CHRYSENE (G5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (G6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (G6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (G6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (G6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (G6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (G6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (G6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CHY 13

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1405	11:34	45	1.003	A BV	28368.	1.391 NG	0.34 y
52	NOT FOUND								
53	NOT FOUND								
54	NOT FOUND								
55	202	1605	13:13	45	1.146	A BB	42024.	2.827 NG	0.69 y
56	240	1848	15:13	56	1.000	A BB	170924.	20.000 NG	4.85
57	202	1643	13:32	56	0.889	A BB	43840.	2.609 NG	0.63 y
58	NOT FOUND								
59	NOT FOUND								
60	228	1852	15:15	56	1.002	A*BB 14200	38454.	3.909 NG	0.95 y
61	149	1852	15:15	56	1.002	A BB	13924.	1.026 NG	0.25 y
62	228	1852	15:15	56	1.002	A VB	19264.	1.984 NG	0.48 y
63	264	2114	17:25	63	1.000	A BB	137048.	20.000 NG	4.85 y
64	NOT FOUND								
65	252	2035	16:45	63	0.963	A BB	28332.	3.596 NG	0.87 y

# SAMPLE DATA PACKAGE

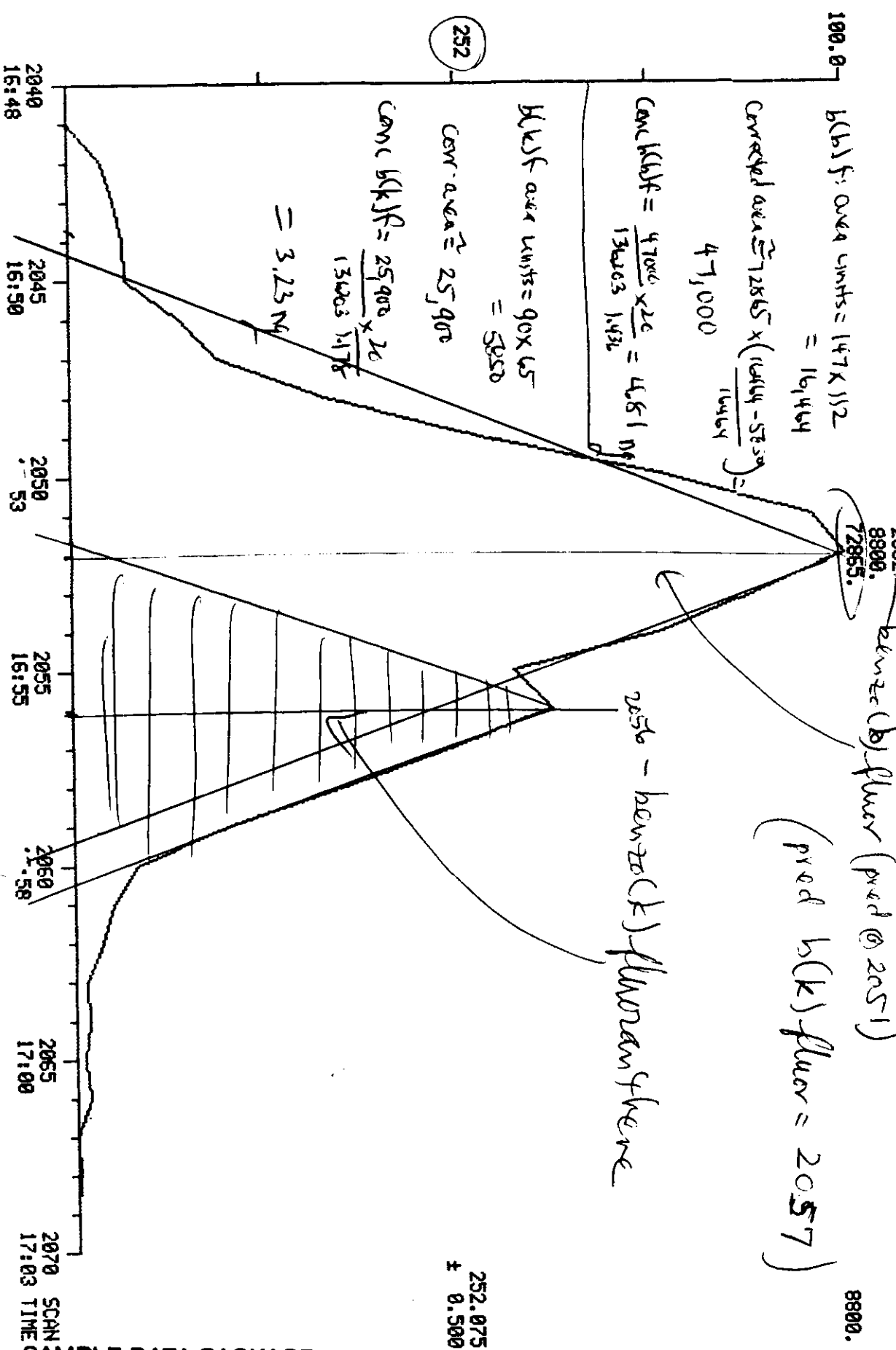
17744 CHY 11

587

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2035	16:45	63	0.963	A BB	28332.	3.596 NG	0.79 y
67	252	2102	17:19	63	0.994	A VB	15080.	2.141 NG	0.52 y
68	276	2412	19:52	63	1.141	A BB	12904.	1.674 NG	0.41 y
69	NOT FOUND								
70	276	2503	20:37	63	1.184	A BB	12060.	1.760 NG	0.43 y
71	112	487	4:01	1	0.727	A BV	302380.	37.696 NG	9.14
72	99	620	5:06	1	0.925	A BV	377685.	37.590 NG	9.11
73	132	638	5:15	1	0.952	A BV	283556.	30.154 NG	7.31
74	152	693	5:42	1	1.034	A BB	188116.	30.146 NG	7.31
75	82	755	6:13	13	0.871	A BB	325917.	31.031 NG	7.52
76	172	1044	8:36	26	0.905	A BV	512132.	31.329 NG	7.60
77	330	1285	10:35	26	1.114	A BB	75332.	26.744 NG	6.48
78	244	1670	13:45	56	0.904	A BV	387008.	35.043 NG	8.50

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:35	1.00	10.000	0.10	1.39	25.00	0.064	1.144	0.06
52	11:53		1.000						
53	11:39		10.000						

MIDWASS CHROMATOGRAM  
 02/11/92 0:33:00  
 SAMPLE: 2UL CC#477484 ID#CHY14 RE Dye/Hz  
 COND.: EXTRACTED 2/4/92 UNOILUTED  
 RANGE: G 1,285 LABEL: N 1, 3.0 QUAN: A 1, 0.1 J 0 BASE: U 20, 3  
 DATA: GR077484C21 #1 SCANS 2040 TO 2070  
 CALL: GR077484C21 #2  
 CS#17744-CHY11 OMR21



52 583 CARBAZOLE  
 53 403 ANTHRACENE (G4#8) <120-12-7>  
 54 426 DI-N-BUTYL PHTHALATE (G4#9) <84-74-2>  
 55 431 FLUORANTHENE (G4#10) <206-44-0>  
 56 \*459 D12-CHRYSENE (IS#5)  
 57 445 PYRENE (G5#3) <129-00-0>  
 58 415 BUTYLBENZYL PHTHALATE (G5#4) <85-68-7>  
 59 423 3,3'-DICHLOROBENZIDINE (G5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (G5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (G5#7) <117-81-7>  
 62 418 CHRYSENE (G5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (G6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (G6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (G6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (G6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (G6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (G6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (G6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CH914

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1417 ✓	11:40	45	1.003	A BV	49724.	4.408 NG	0.85 YES
52	NOT FOUND								
53	178	1426 ✓	11:45	45	1.009	A VV	9048.	1.033 NG	0.20 YES
54	NOT FOUND								
55	202	1617	13:19	45	1.144	A BV	102051.	10.641 NG	2.04 YES
56	240	1861	15:20	56	1.000	A BB	193728.	20.000 NG	3.84 YES
57	202	1655	13:38	56	0.889	A BB	82547.	6.041 NG	1.16 YES
58	149	1764	14:32	56	0.948	A BB	8308.	1.690 NG	0.32 YES
59	NOT FOUND								
60	228	1860	15:19	56	0.999	A*BV	60174 166888.	5.639 9.457 NG	1.82 YES
61	149	1863	15:21	56	1.001	A BB	53227.	4.728 NG	0.91 YES
62	228	1860	15:19	56	0.999	A*BV	36363 166888.	4.512 12.513 NG	2.40 YES
63	264	2136	17:35	63	1.000	A BB	136203.	20.000 NG	3.84 YES
64	NOT FOUND								
65	252	2052	16:54	63	0.961	A BV	47000 80828.	4.81 = 360 ug 9.264 NG	1.59 YES

SAMPLE DATA PACKAGE

17744 CHY II

653

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2056	16:54	63	0.961	A BV	25900 80828.	3.23 = 240 ug 10.079 NG	1.93 YES
67	252	2123	17:29	63	0.994	A VV	31180.	4.863 NG	0.93 YES
68	276	2444	20:08	63	1.144	A BB	19004.	2.965 NG	0.57 YES
69	278	2451	20:11	63	1.147	A BB	6192.	1.044 NG	0.20 YES
70	276	2541	20:56	63	1.190	A*BB	15940.	2.588 NG	0.50 YES
71	112	500	4:07	1	0.735	A BV	339320.	51.877 NG	9.96
72	99	630	5:11	1	0.926	A BV	385783.	49.439 NG	9.49
73	132	648	5:20	1	0.953	A BB	255616.	40.251 NG	7.73
74	152	703	5:47	1	1.034	A BB	138600.	29.872 NG	5.73
75	82	765	6:18	13	0.872	A BB	310960.	37.287 NG	7.16
76	172	1054	8:41	26	0.905	A BB	351540.	31.838 NG	6.11
77	330	1296	10:40	26	1.112	A BB	77132.	29.455 NG	5.65
78	244	1682	13:51	56	0.904	A BB	328490.	39.468 NG	7.58

63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (G6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (G6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (G6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (G6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (G6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (G6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (G6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

STD  
 2/11

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1418	11:41	45	1.003	A BV	302272.	24.822 NG	1.33
52	167	1456	11:59	45	1.030	A BB	170508.	25.265 NG	1.35
53	178	1426	11:45	45	1.008	A VV	260924.	25.216 NG	1.35
54	149	1513	12:28	45	1.070	A BV	362744.	24.845 NG	1.33
55	202	1618	13:19	45	1.144	A BV	225092.	25.894 NG	1.38
56	240	1861	15:20	56	1.000	A BB	90116.	20.000 NG	1.07
57	202	1656	13:38	56	0.890	A BV	231136.	26.345 NG	1.41
58	149	1765	14:32	56	0.948	A BV	103288.	25.302 NG	1.35
59	252	1855	15:17	56	0.997	A BB	7964.	24.293 NG	1.30
60	228	1860	15:19	56	0.999	A BV	128244.	26.041 NG	1.39
61	149	1864	15:21	56	1.002	A BB	169928.	24.308 NG	1.30
62	228	1864	15:22	56	1.003	A VB	120264.	24.800 NG	1.33
63	264	2133	17:34	63	1.000	A BB	52960.	20.000 NG	1.07
64	149	1976	16:16	63	0.926	A BB	171824.	23.529 NG	1.26
65	252	2049	16:52	63	0.961	A BV	82496.	25.885 NG	1.38

IS S

IS 6

b(b) f

# SAMPLE DATA PACKAGE

17744 CHY 11

1749

Quantitation Report File: HQ920211B21

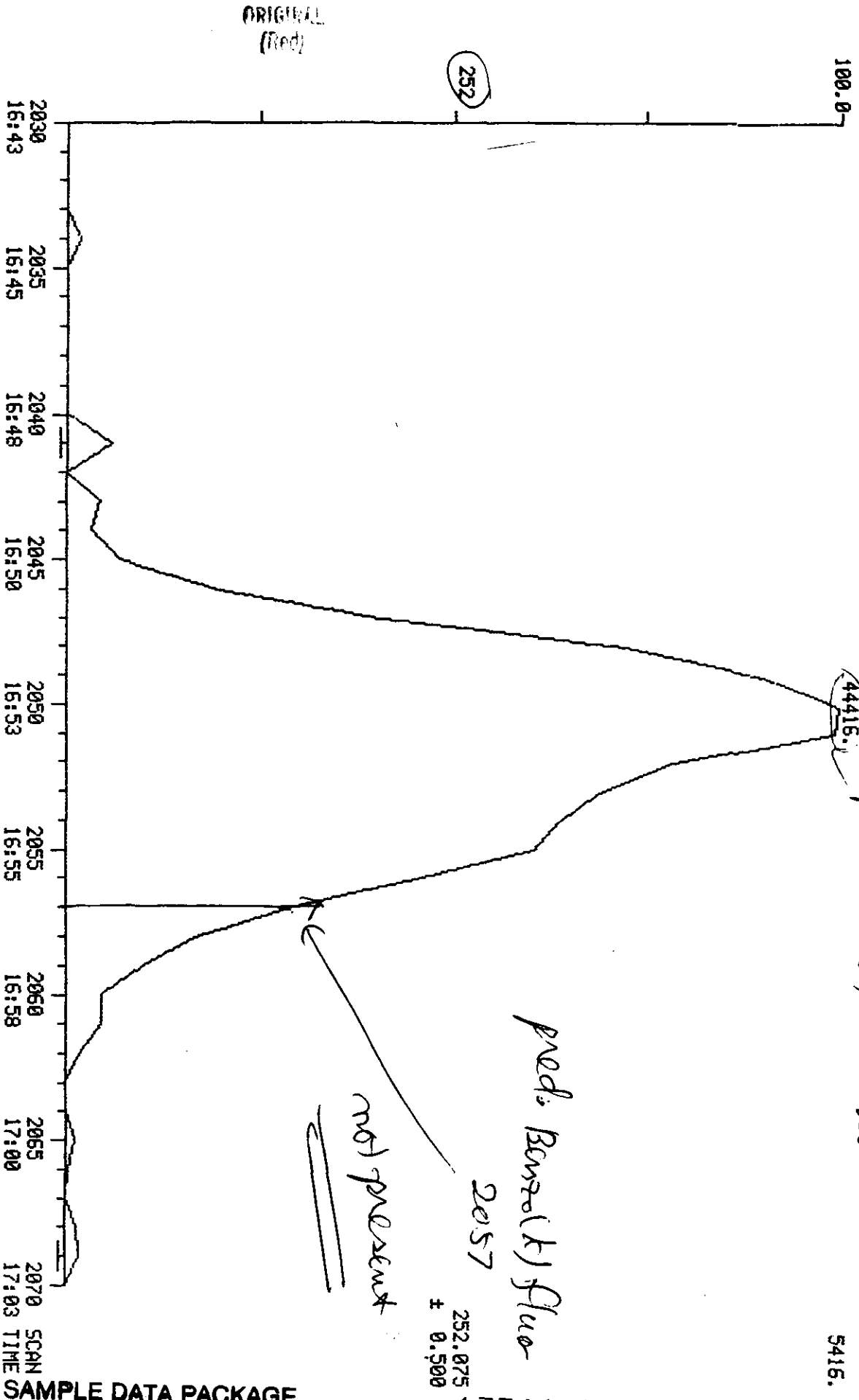
No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2055	16:55	63	0.963	A VB	80580.	12.167 NG	0.65
67	252	2120	17:27	63	0.994	A BV	63072.	23.192 NG	1.24
68	276	2438	20:05	63	1.143	A BB	65380.	24.477 NG	1.31
69	278	2445	20:08	63	1.146	A BB	53460.	25.574 NG	1.37
70	276	2532	20:51	63	1.187	A BB	56044.	24.345 NG	1.30
71	112	495	4:05	1	0.728	A BV	170553.	24.080 NG	1.29
72	99	628	5:10	1	0.924	A BB	207764.	23.697 NG	1.27
73	132	647	5:20	1	0.951	A BB	164984.	24.038 NG	1.28
74	152	703	5:47	1	1.034	A BB	121528.	24.521 NG	1.31
75	82	765	6:18	13	0.871	A VB	226532.	23.483 NG	1.26
76	172	1055	8:41	26	0.905	A BB	285048.	25.934 NG	1.39
77	330	1296	10:40	26	1.111	A BB	33340.	28.782 NG	1.54
78	244	1682	13:51	56	0.904	A BV	133732.	26.487 NG	1.42

b(k) f

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:41	1.00	10.000	0.10	24.82	25.00	1.142	1.150	0.99
52	11:59	1.00	1.000	1.03	25.27	25.00	0.644	0.637	1.01
53	11:45	1.00	10.000	0.10	25.22	25.00	0.986	0.977	1.01
54	12:28	1.00	10.000	0.11	24.85	25.00	1.371	1.379	0.99
55	13:19	1.00	10.000	0.11	25.89	25.00	0.850	0.821	1.04
56	15:20	1.00	10.000	0.10	20.00	20.00	1.000	1.000	1.00
57	13:39	1.00	10.000	0.09	26.34	25.00	2.052	1.947	1.05



MIDMASS CHROMATOGRAM  
 02/12/92 5:20:00  
 SAMPLE: 2UL CC#477485 ID#CHY15 R5  
 CONDS.: EXTRACTED 2/4/92 1: 25 DILUTION  
 RANGE: G 1,2855 LABEL: H 1, 3.0 QUANT: A 1, 0.13 0  
 DATA: GR077485A21 #1  
 CALL: GR077485A21 #2  
 CS#17744-CHY11  
 OM21  
 SCANS 2030 TO 2070  
 BASE: U 20, 3



SAMPLE DATA PACKAGE

17744 CHY 11

58 415 BUTYLBENZYL PHTHALATE (Q5#4) <85-68-7>  
 59 423 3,3'-DICHLOROBENZIDINE (Q5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (Q5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (Q5#7) <117-81-7>  
 62 418 CHRYSENE (Q5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (Q6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (Q6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (Q6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (Q6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (Q6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (Q6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (Q6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CHY 15

	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1417	11:40	45	1.002	A VV	34388.	2.902 NG	1.37 yes
52	167	1456	11:59	45	1.030	A BV	18256.	2.731 NG	1.29 NO
53	178	1427	11:45	45	1.009	A*VB	13574.	1.327 NG	0.63 NO
54	NOT FOUND								
55	202	1618	13:19	45	1.144	A BB	62086.	7.035 NG	3.32 yes
56	240	1861	15:20	56	1.000	A BV	122564.	20.000 NG	9.45 155
57	202	1657	13:39	56	0.890	A BB	60743.	4.831 NG	2.28 yes
58	NOT FOUND								
59	252	1858	15:18	56	0.998	A BV	3856.	8.900 NG	4.20 NO
60	228	1865	15:21	56	1.002	A*BB	20616 49740 2.956	7.129 NG	3.37 yes
61	149	1864	15:21	56	1.002	A BV	11926.	1.290 NG	0.61 yes
62	228	1865	15:21	56	1.002	A*BB	24652 49740 3.767	7.602 NG	3.59 yes
63	264	2133	17:34	63	1.000	A BV	91208.	20.000 NG	9.45 156
64	NOT FOUND								
65	252	2050	16:53	63	0.961	A BB	44416. 7.905	7.816 NG	3.69 yes

= 7800 wj/kg

1604  
72242

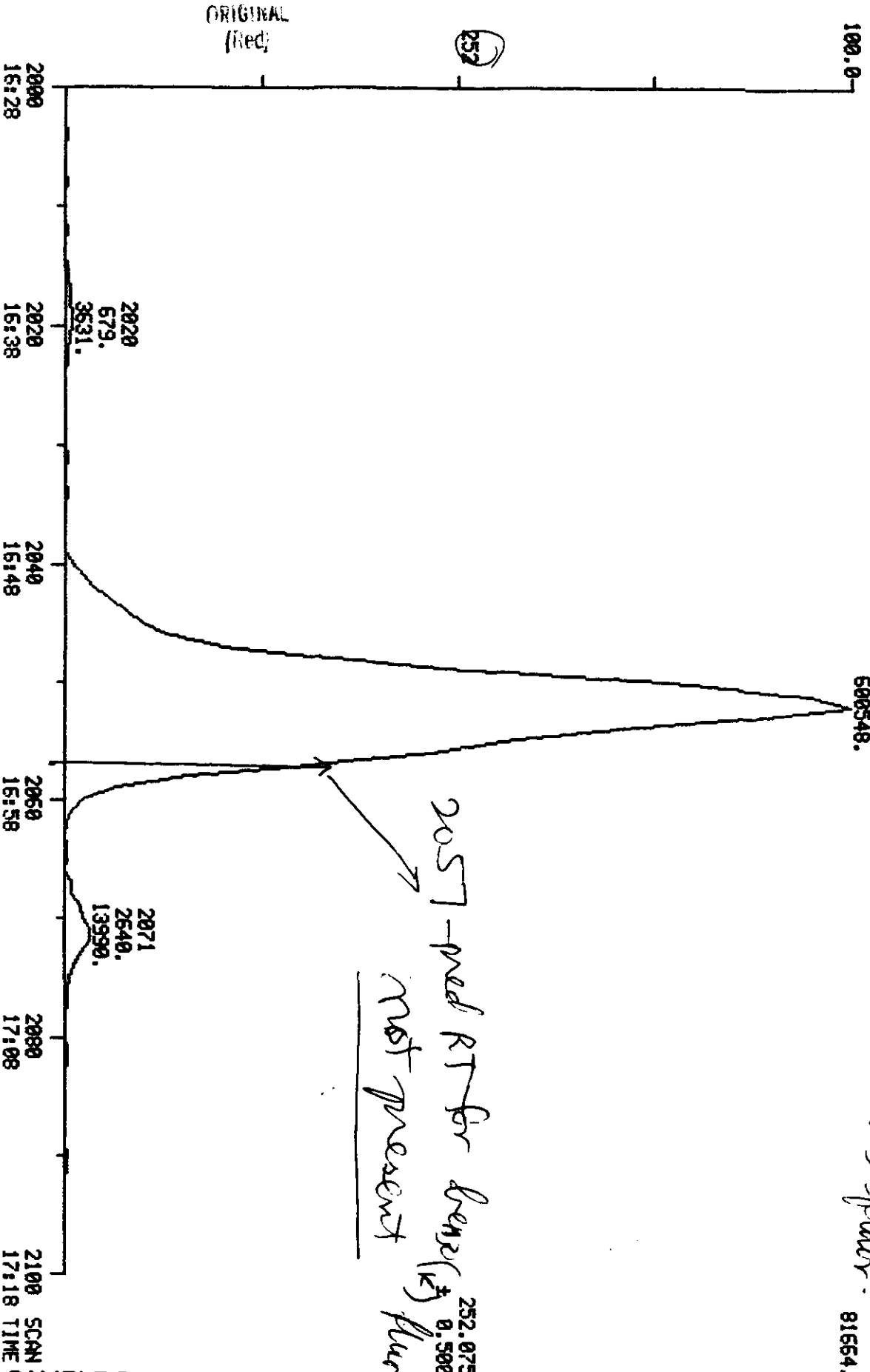
# SAMPLE DATA PACKAGE

17744 CHY 11

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2050	16:53	63	0.961	A BB	44416. 7.905	7.816 NG	3.78 yes
67	252	2120	17:27	63	0.994	A VB	15892.	3.658 NG	1.73 yes
68	276	2441	20:06	63	1.144	A BV	10664.	2.368 NG	1.12 yes
69	NOT FOUND								
70	276	2532	20:51	63	1.187	A*BV	8184.	2.120 NG	1.00 yes
71	NOT FOUND								
72	NOT FOUND								
73	NOT FOUND								
74	NOT FOUND								
75	NOT FOUND								
76	NOT FOUND								
77	330	1297	10:41	26	1.114	A BB	2328.	1.925 NG	0.91
78	244	1682	13:51	56	0.904	A BB	12660.	1.740 NG	0.82

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:41	1.00	10.000	0.10	2.90	25.00	0.133	1.142	0.12
52	11:59	1.00	1.000	1.03	2.73	25.00	0.070	0.644	0.11
53	11:45	1.00	10.000	0.10	1.33	25.00	0.052	0.986	0.05
54	12:28		10.000						

MIDWASS CHROMATOGRAM  
 02/10/92 12:58:00  
 SAMPLE: 2UL CC#477489 ID#CHY18.82  
 CONDS.: EXTRACTED 2/4/92 UNOILUTED  
 RANGE: G 1,2855 LABEL: N 1, 4.0 QUAN: A 2, 1.0 J 0  
 DATA: GR077489C21 #1 SCANS 2000 TO 2100  
 CALL: GR077489C21 #2  
 CS#17744-CHY11  
 QM#21



SAMPLE DATA PACKAGE

17744 CHY 11

57 400 BENZO(A)PYRENE (G6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (G6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (G6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (G6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CHY18 initd

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1416	11:40	45	1.003	A BV	341736.	27.221 NG	2.94 yes
52	167	1453	11:58	45	1.029	A VB	32036.	3.924 NG	0.42 yes
53	178	1425	11:44	45	1.009	A VV	32672.	2.962 NG	0.32 yes
54	NOT FOUND								
55	202	1618	13:19	45	1.146	A BB	795781.	80.740 NG	8.73 yes E
56	240	1861	15:20	56	1.000	A BB	189320.	20.000 NG	2.16
57	202	1656	13:38	56	0.890	A BB	732198.	58.935 NG	6.37 yes
58	NOT FOUND								
59	NOT FOUND								
60	228	1859	15:19	56	0.999	A BV	348068.	36.564 NG	3.95 yes
61	149	1862	15:20	56	1.001	A BB	13272.	1.691 NG	0.18 no
62	228	1865	15:21	56	1.002	A VV	308040.	34.900 NG	3.77 yes
63	264	2134	17:34	63	1.000	A BB	98796.	20.000 NG	2.16
64	NOT FOUND								
65	252	2052	16:54	63	0.962	A BV	625398. 91.905 NG	91.905 NG	9.94 yes

4100 v9/kg

# SAMPLE DATA PACKAGE

17744 CHY 11

9030

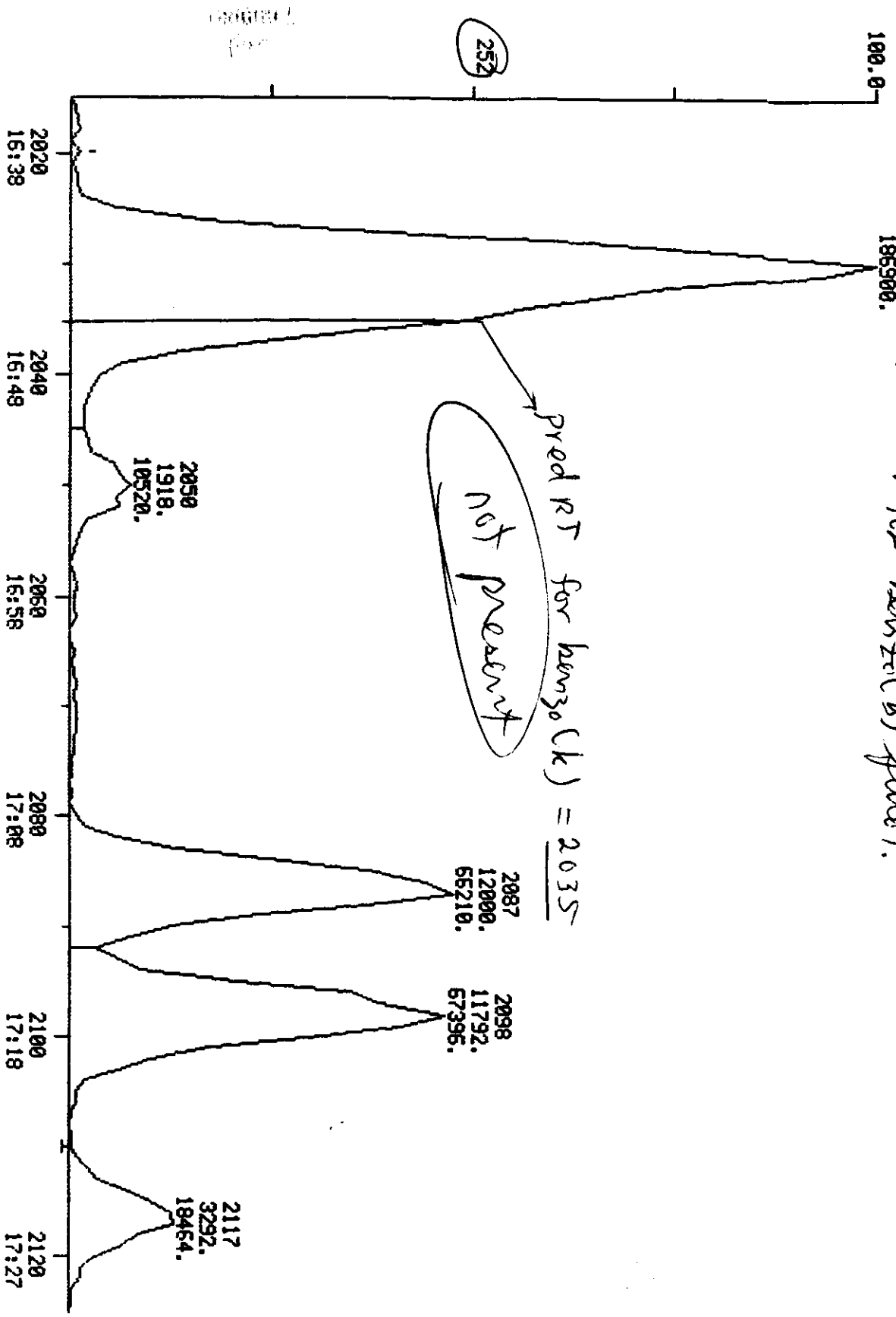
No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2052	16:54	63	0.962	A BV	625398.	91.905 NG	12.87 yes
67	252	2122	17:28	63	0.994	A VB	137591.	26.938 NG	2.91 yes
68	276	2442	20:07	63	1.144	A BB	123538.	21.525 NG	2.33 yes
69	NOT FOUND								
70	276	2536	20:53	63	1.188	A BB	101238.	20.881 NG	2.26 yes
71	112	499	4:07	1	0.734	A BB	166604.	36.171 NG	3.91
72	99	629	5:11	1	0.925	A BV	241168.	44.336 NG	4.79
73	132	648	5:20	1	0.953	A BV	167116.	31.390 NG	3.39
74	152	703	5:47	1	1.034	A BB	116704.	31.701 NG	3.43
75	82	764	6:17	13	0.871	A BV	214856.	39.569 NG	4.28
76	172	1053	8:40	26	0.905	A BB	371204.	32.096 NG	3.47
77	330	1295	10:40	26	1.113	A BB	27776.	13.530 NG	1.46
78	244	1681	13:51	56	0.903	A BB	337416.	39.662 NG	4.29

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:42	1.00	10.000	0.10	27.22	25.00	1.251	1.149	1.09
52	12:00	1.00	1.000	1.03	3.92	25.00	0.117	0.747	0.16
53	11:47	1.00	10.000	0.10	2.96	25.00	0.120	1.010	0.12
54	12:29		10.000						
55	13:21	1.00	10.000	0.11	80.74	25.00	2.914	0.902	3.23
56	15:21	1.00	10.000	0.10	20.00	20.00	1.000	1.000	1.00
57	13:40	1.00	10.000	0.09	58.94	25.00	3.094	1.312	2.36
58	14:34		10.000						
59	15:18		20.000						
60	15:21	1.00	10.000	0.10	36.56	25.00	1.471	1.006	1.46
61	15:22	1.00	10.000	0.10	1.69	25.00	0.056	0.829	0.07
62	15:23	1.00	10.000	0.10	34.90	25.00	1.302	0.932	1.40
63	17:36	1.00	10.000	0.10	20.00	20.00	1.000	1.000	1.00
64	16:17		10.000						
65	16:55	1.00	10.000	0.10	91.905	25.00			

no RF-

MIDWASS CHROMATOGRAM  
 02/12/92 20:55:00  
 SAMPLE: 2UL CC#477489 ID#CHY18 RE  
 COND.: EXTRACTED 2/4/92 1: 2 DILUTION  
 RANGE: G 1.2855 LABEL: N 1, 3.0 QUAN: A 1, 0.3 J 0 BASE: U 20, 3  
 DATA: GDJ77489R21 #1  
 CALL: GDJ77489R21 #2  
 CS#17744-CHY11  
 OMA21  
 SCANS 2015 TO 2125

25216. → pred. RT for benz(c)b fluor.  
 185900.



SAMPLE DATA PACKAGE

252.075  
 ± 0.500  
 17744 CHY 11

25216.

58 413 BUTYLBENZYL PHTHALATE (Q5#4) <85-68-7>  
 59 423 3,3'-DICHLOROBENZIDINE (Q5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (Q5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (Q5#7) <117-81-7>  
 62 418 CHRYSENE (Q5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (Q6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (Q6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (Q6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (Q6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (Q6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (Q6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (Q6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CN418DL

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1401	11:32	45	1.003	A BV	170436.	14.871 NG	2.43 ycs
52	167	1438	11:51	45	1.029	A BB	23228.	2.845 NG	0.47 ycs
53	178	1409	11:36	45	1.009	A VV	25788.	2.586 NG	0.42 ycs
54	NOT FOUND								
55	202	1601	13:11	45	1.146	A BV	358208.	42.890 NG	7.02 ycs
56	240	1844	15:11	56	1.000	A BB	85720.	20.000 NG	3.27 ycs
57	202	1639	13:30	56	0.889	A BB	304797.	36.168 NG	5.92 ycs
58	NOT FOUND								
59	NOT FOUND								
60	228	1842	15:10	56	0.999	A BV	91680.	18.580 NG	3.04 ycs
61	149	1848	15:13	56	1.002	A BB	15852.	2.328 NG	0.38 ycs
62	228	1848	15:13	56	1.002	A VV	109664.	22.520 NG	3.68 ycs
63	264	2110	17:23	63	1.000	A BV	64684.	20.000 NG	3.27 ycs
64	NOT FOUND								
65	252	2030	16:43	63	0.962	A BV	1460187097A 47.913	50.811 NG	8.23 ycs

= 4500 ycs/kg

969

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2098	17:17	63	0.994	A VV	68668.	20.655 NG	3.38 ycs
67	252	2098	17:17	63	0.994	A VV	68668.	20.655 NG	3.38 ycs
68	276	2408	19:50	63	1.141	A BB	64357 79212.	19.073 21.777 NG	3.56 ycs
69	278	2414	19:53	63	1.144	A VB	19089 26772.	6.468 9.055 NG	1.48 ycs
70	276	2501	20:36	63	1.185	A BB	65400 72032.	20.221 22.275 NG	3.64 ycs
71	112	483	3:59	1	0.725	A BB	116756.	25.493 NG	4.17
72	99	615	5:04	1	0.923	A BB	148084.	25.814 NG	4.22
73	132	633	5:13	1	0.950	A BB	110520.	20.585 NG	3.37
74	152	688	5:40	1	1.033	A BB	60392.	16.950 NG	2.77
75	82	750	6:11	13	0.869	A BB	126120.	20.927 NG	3.42
76	172	1039	8:33	26	0.904	A BB	173808.	18.830 NG	3.08
77	330	1280	10:32	26	1.114	A BB	27244.	17.129 NG	2.80
78	244	1665	13:43	56	0.903	A BB	126276.	22.799 NG	3.73

Opul

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:35	1.00	10.000	0.10	14.87	25.00	0.681	1.144	0.59
52	11:53	1.00	1.000	1.03	2.84	25.00	0.093	0.815	0.11
53	11:39	1.00	10.000	0.10	2.59	25.00	0.103	0.996	0.10
54	12:22		10.000						
55	13:14	1.00	10.000	0.11	42.89	25.00	1.431	0.834	1.72

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

EPA SAMPLE NO.

Lab Name: COMPUCHEM RTP

Contract: 68D10083

CHY18

CHY18DL

Lab Code: COMPU Case No.: 17744

SAS No.: 6579HQ

SDG No.: CHY11

3 No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477489

: 477489

Sample wt/vol: 30.4 (g/mL) G

Lab File ID: GR077489C21

GDJ77489A21

Level: (low/med) LOW

Date Received: 01/29/92

: 01/29/92

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 02/04/92

1: 02/04/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 02/10/92

: 02/12/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

or: 2.0

GPC Cleanup: (Y/N) Y pH: 6.1

CAS NO.

COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Q

FS:  
3/KG

Q

108-95-2-----	Phenol	450	U	900	U
111-44-4-----	bis(2-Chloroethyl) Ether	450	U	900	U
95-57-8-----	2-Chlorophenol	450	U	900	U
541-73-1-----	1,3-Dichlorobenzene	450	U	900	U
106-46-7-----	1,4-Dichlorobenzene	450	U	900	U
95-50-1-----	1,2-Dichlorobenzene	450	U	900	U
95-48-7-----	2-Methylphenol	450	U	900	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	450	U	900	U
106-44-5-----	4-Methylphenol	450	U	900	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	450	U	900	U
67-72-1-----	Hexachloroethane	450	U	900	U
98-95-3-----	Nitrobenzene	450	U	900	U
78-59-1-----	Isophorone	450	U	900	U
88-75-5-----	2-Nitrophenol	450	U	900	U
105-67-9-----	2,4-Dimethylphenol	450	U	900	U
111-91-1-----	bis(2-Chloroethoxy)Methane	450	U	900	U
120-83-2-----	2,4-Dichlorophenol	450	U	900	U
120-82-1-----	1,2,4-Trichlorobenzene	450	U	900	U
91-20-3-----	Naphthalene	450	U	900	U
106-47-8-----	4-Chloroaniline	450	U	900	U
87-68-3-----	Hexachlorobutadiene	450	U	900	U
59-50-7-----	4-Chloro-3-Methylphenol	450	U	900	U
91-57-6-----	2-Methylnaphthalene	450	U	900	U
77-47-4-----	Hexachlorocyclopentadiene	450	U	900	U
88-06-2-----	2,4,6-Trichlorophenol	450	U	900	U
95-95-4-----	2,4,5-Trichlorophenol	1100	U	2200	U
91-58-7-----	2-Chloronaphthalene	450	U	900	U
88-74-4-----	2-Nitroaniline	1100	U	2200	U
131-11-3-----	Dimethyl Phthalate	450	U	900	U
208-96-8-----	Acenaphthylene	450	U	900	U
606-20-2-----	2,6-Dinitrotoluene	450	U	900	U
99-09-2-----	3-Nitroaniline	1100	U	2200	U
83-32-9-----	Acenaphthene	89	J	110	DJ

FORM I SV-1

SAMPLE DATA PACKAGE

17744 CHY 11

all surr received in ltrh 3/90  
(except TRP in 895 initial anal) 960

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CHY18

CHY18DL

Lab Code: COMPU

Case No.: 17744

SAS No.: 6579HQ

SDG No.: CHY11

XG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477489

ID: 477489

Sample wt/vol: 30.4 (g/mL) G

Lab File ID: GR077489C21

GDJ77489A21

Level: (low/med) LOW

Date Received: 01/29/92

ID: 01/29/92

% Moisture: 28 decanted: (Y/N) N

Date Extracted: 02/04/92

ID: 02/04/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 02/10/92

ID: 02/12/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

ID: 2.0

GPC Cleanup: (Y/N) Y

pH: 6.1

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

ITS:  
UG/KG

CAS NO.	COMPOUND	(ug/L or ug/Kg) UG/KG	Q	ITS: UG/KG	Q
51-28-5-----	2,4-Dinitrophenol	1100	U	2200	U
100-02-7-----	4-Nitrophenol	1100	U	2200	U
132-64-9-----	Dibenzofuran	450	U	900	U
121-14-2-----	2,4-Dinitrotoluene	450	U	900	U
84-66-2-----	Diethylphthalate	450	U	390	DJ
7005-72-3-----	4-Chlorophenyl-phenylether	450	U	900	U
86-73-7-----	Fluorene	66	J	92	DJ
100-01-6-----	4-Nitroaniline	1100	U	2200	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1100	U	2200	U
86-30-6-----	N-Nitrosodiphenylamine (1)	450	U	900	U
101-55-3-----	4-Bromophenyl-phenylether	450	U	900	U
118-74-1-----	Hexachlorobenzene	450	U	900	U
87-86-5-----	Pentachlorophenol	1100	U	2200	U
85-01-8-----	Phenanthrene	1200	U	1400	D
120-12-7-----	Anthracene	140	J	240	DJ
86-74-8-----	Carbazole	180	J	260	DJ
84-74-2-----	Di-n-Butylphthalate	450	U	900	U
206-44-0-----	Fluoranthene	3700	E	3900	D
129-00-0-----	Pyrene	2700	U	3300	D
85-68-7-----	Butylbenzylphthalate	450	U	900	U
91-94-1-----	3,3'-Dichlorobenzidine	450	U	900	U
56-55-3-----	Benzo(a)Anthracene	1700	U	1700	D
218-01-9-----	Chrysene	1600	U	2100	D
117-81-7-----	bis(2-Ethylhexyl)Phthalate	450	U	900	U
117-84-0-----	Di-n-Octyl Phthalate	450	U	900	U
205-99-2-----	Benzo(b)Fluoranthene	450	U	900	U
207-08-9-----	Benzo(k)Fluoranthene	450	U	900	U
50-32-8-----	Benzo(a)Pyrene	1200	U	1900	D
193-39-5-----	Indeno(1,2,3-cd)Pyrene	980	U	1700	D
53-70-3-----	Dibenz(a,h)Anthracene	350	J	590	DJ
191-24-2-----	Benzo(g,h,i)Perylene	950	U	1800	D

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

SAMPLE DATA PACKAGE

17744 CHY 11

896

3/90

96



MIDMASS CHROMATOGRAM  
 02/12/92 4:05:00  
 SAMPLE: 2UL CC#477490 ID#CHY19 RE  
 COND5.: EXTRACTED 2/4/92 1: 2 DILUTION  
 RANGE: G 1,2855 LABEL: N 1, 3.0 QUAM: A 1, 0.1 J 0 BASE: U 20, 3  
 DATA: GR077490R21 #1 SCANS 2030 TO 2070  
 CALL: GR077490R21 #2  
 CS#17744-CHY11 QMR21

Conv area for h(b)f:

$$152616 \times \left( \frac{12035 - 3360}{12035} \right) = 109,000$$

Conv conc:

$$\frac{109,000}{70740} \times \frac{20}{1.296} = 24.6 \text{ ng}$$

Conv can for b(k)f:

$$\frac{41656}{70740} \times \frac{20}{1.217} = 9.67 \text{ ng}$$

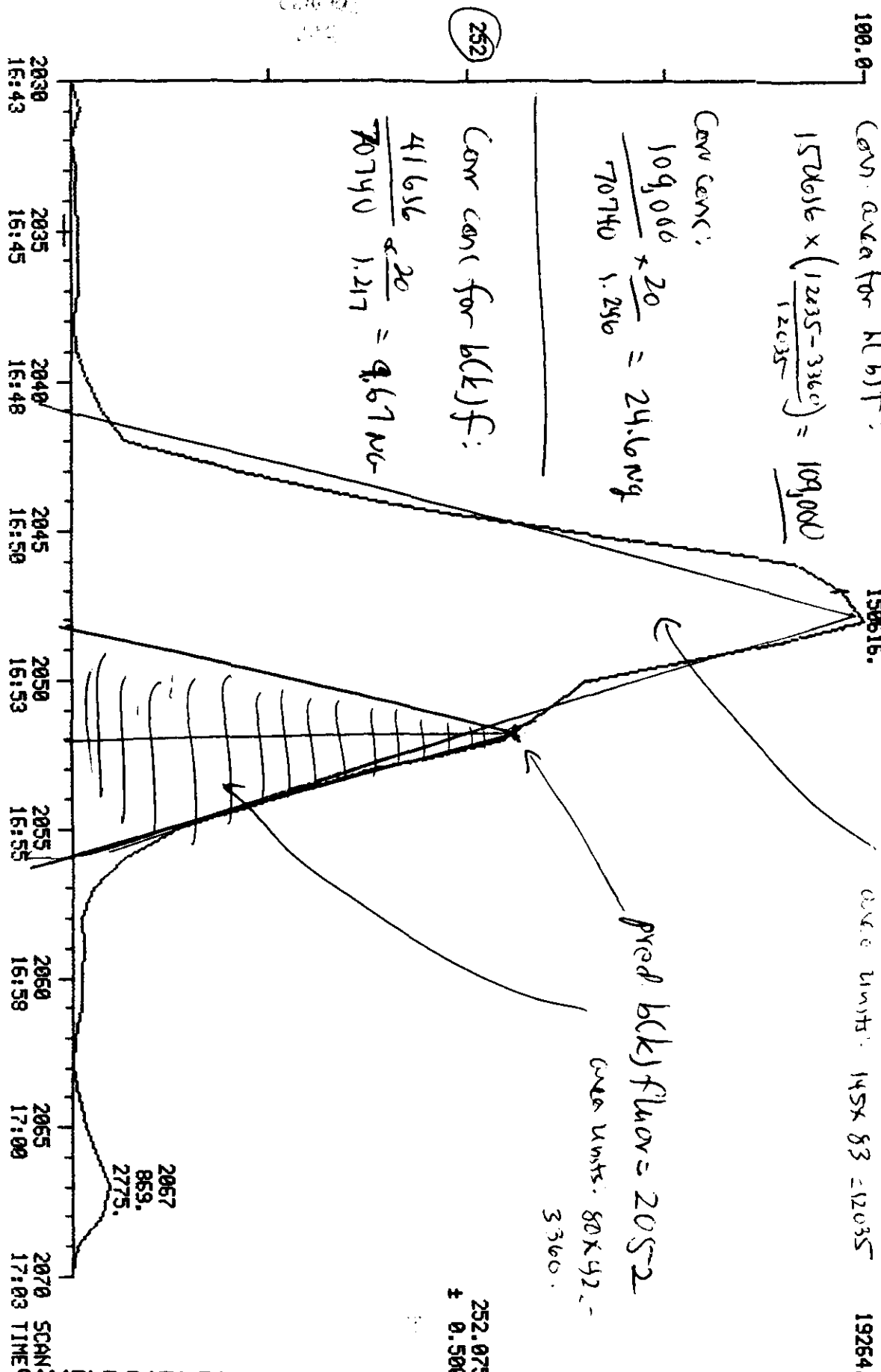
pred b(k) fluor = 2046

area units: 145x 83 = 12035 19264.

pred b(k) fluor = 2052

area units: 80x42 = 3360.

252.075  
 ± 0.500



55 431 FLUORANTHENE (G4#10) <206-44-0>  
 56 \*459 D12-CHRYSENE (IS#3)  
 57 445 PYRENE (G5#3) <129-00-0>  
 58 415 BUTYLBENZYL PHTHALATE (G5#4) <85-68-7>  
 59 423 3,3'-DICHLOROBENZIDINE (G5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (G5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (G5#7) <117-81-7>  
 62 418 CHRYSENE (G5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (G6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (G6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (G6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (G6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (G6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (G6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (G6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 /8 #496 D14-TERPHENYL (SS#8)

CHY 19

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1413 ✓	11:38	45	1.003	A BV	219332.	26.919 NG	5.07 yes
52	167	1450	11:56	45	1.029	A BB	17400.	3.786 NG	0.71 yes
53	178	1422 ✓	11:43	45	1.009	A VV	36232.	5.151 NG	0.97 yes
54	NOT FOUND								
55	202	1614 ✓	13:17	45	1.145	A BV	284737.	46.929 NG	8.84 yes
56	240	1857	15:18	56	1.000	A BB	96780.	20.000 NG	3.77
57	202	1652 ✓	13:36	56	0.890	A BB	228605.	23.024 NG	4.34 yes
58	NOT FOUND								
59	NOT FOUND								
60	228	1856 ✓	15:17	56	0.999	A BV	87168.	15.823 NG	2.98 yes
61	NOT FOUND								
62	228	1852 ✓	15:20	56	1.003	A VV	91120.	17.637 NG	3.32 yes
63	264	2130	17:32	63	1.000	A BV	70740.	20.000 NG	3.77
64	NOT FOUND								
65	252	2048	16:52	63	0.962	A BV	109,000 109,000	24.6 NG 24.6 NG	6.53 yes

155

156

x86 = 2100 ug/kg

164

# SAMPLE DATA PACKAGE

17744 CHY 11

1044-1442

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2048	16:52	63	0.962	A BV	109,000	24.6 NG	6.68 yes
67	252	2118	17:27	63	0.994	A VV	60634.	17.993 NG	3.39 yes
68	276	2438	20:05	63	1.145	A BB	43772.	12.531 NG	2.36 yes
69	278	2448	20:10	63	1.149	A BV	13217.	4.627 NG	0.87 yes
70	276	2533	20:52	63	1.189	A VB	34592.	11.552 NG	2.18 yes
71	112	492	4:03	1	0.729	A BB	87308.	19.269 NG	3.63
72	99	625	5:09	1	0.926	A BV	133016.	24.099 NG	4.54
73	132	643	5:18	1	0.953	A BB	71104.	16.223 NG	3.06
74	152	698	5:45	1	1.034	A BB	52036.	16.118 NG	3.04
75	82	760	6:16	13	0.871	A BB	108228.	19.237 NG	3.62
76	172	1050	8:39	26	0.904	A BB	140344.	17.799 NG	3.35
77	330	1293	10:39	26	1.114	A BB	7336.	7.955 NG	1.50
78	244	1678	13:49	56	0.904	A BB	95012.	16.539 NG	3.12

9.67 x86 = 830 ug/kg

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:41	1.00	10.000	0.10	26.92	25.00	1.230	1.142	1.08
52	11:59	1.00	1.000	1.03	3.79	25.00	0.098	0.644	0.15

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

EPA SAMPLE NO.

CHY20

CHY20DL

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477491

ID: 477491

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GRD77491C21

: GRD77491A21

Level: (low/med) LOW Date Received: 01/29/92

ad: 01/29/92

% Moisture: 35 decanted: (Y/N) N Date Extracted: 02/04/92

ted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/11/92

ed: 02/13/92

Injection Volume: 2.0 (uL) Dilution Factor: 15.0

stor: 30.0

GPC Cleanup: (Y/N) Y pH: 6.7

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

NITS:  
UG/KG

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q	Q
51-28-5	2,4-Dinitrophenol	18000	U	36000
100-02-7	4-Nitrophenol	18000	U	36000
132-64-9	Dibenzofuran	2900	J ✓	3100
121-14-2	2,4-Dinitrotoluene	7500	U	15000
84-66-2	Diethylphthalate	7500	U ✓	3000
7005-72-3	4-Chlorophenyl-phenylether	7500	U	15000
86-73-7	Fluorene	4100	J ✓	4900
100-01-6	4-Nitroaniline	18000	U	36000
534-52-1	4,6-Dinitro-2-Methylphenol	18000	U	36000
86-30-6	N-Nitrosodiphenylamine (1)	7500	U	15000
101-55-3	4-Bromophenyl-phenylether	7500	U	15000
118-74-1	Hexachlorobenzene	7500	U	15000
87-86-5	Pentachlorophenol	18000	U	36000
85-01-8	Phenanthrene	75000	E	52000
120-12-7	Anthracene	14000	✓	13000
86-74-8	Carbazole	11000	✓	7900
84-74-2	Di-n-Butylphthalate	7500	U	15000
206-44-0	Fluoranthene	110000	E	72000
129-00-0	Pyrene	51000	✓	54000
85-68-7	Butylbenzylphthalate	7500	U	15000
91-94-1	3,3'-Dichlorobenzidine	7500	U	15000
56-55-3	Benzo(a)Anthracene	33000	✓	29000
218-01-9	Chrysene	32000	✓	27000
117-81-7	bis(2-Ethylhexyl)Phthalate	1600	33 ✓	2600
117-84-0	Di-n-Octyl Phthalate	7500	U	15000
205-99-2	Benzo(b)Fluoranthene	46000	✓	36,000
207-08-9	Benzo(k)Fluoranthene	15000	✓	9700
50-32-8	Benzo(a)Pyrene	30000	✓	24000
193-39-5	Indeno(1,2,3-cd)Pyrene	11000	✓	15000
53-70-3	Dibenz(a,h)Anthracene	5300	J ✓	6000
191-24-2	Benzo(g,h,i)Perylene	7600	✓	13000

(1) - Cannot be separated from Diphenylamine

FORM I SV-2 *surrogate recoveries 3/90*

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1109

1182

1B  
ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

EPA SAMPLE NO.

P Contract: 68D10083

CHY20

CHY20DL

ase No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

SDG No.: CHY11

OIL Lab Sample ID: 477491

ID: 477491

30.5 (g/mL) G Lab File ID: GRD77491C21

: GDJ77491A21

LOW Date Received: 01/29/92

ed: 01/29/92

ecanted: (Y/N) N Date Extracted: 02/04/92

ted: 02/04/92

Volume: 500.0 (uL) Date Analyzed: 02/11/92

ed: 02/13/92

2.0(uL) Dilution Factor: 15.0

ctor: 30.0

pH: 6.7

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

NITS:  
UG/KG Q

COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q		NITS: UG/KG	Q
-Phenol	7500	U		15000	U
-bis(2-Chloroethyl) Ether	7500	U		15000	U
--2-Chlorophenol	7500	U		15000	U
-1,3-Dichlorobenzene	7500	U		15000	U
-1,4-Dichlorobenzene	7500	U		15000	U
--1,2-Dichlorobenzene	7500	U		15000	U
--2-Methylphenol	7500	U		15000	U
-2,2'-Oxybis(1-Chloropropane)	7500	U		15000	U
-4-Methylphenol	7500	U		15000	U
--N-Nitroso-Di-n-Propylamine	7500	U		15000	U
-Hexachloroethane	7500	U		15000	U
-Nitrobenzene	7500	U		15000	U
-sophorone	7500	U		15000	U
--2-Nitrophenol	7500	U		15000	U
-2,4-Dimethylphenol	7500	U		15000	U
-bis(2-Chloroethoxy)Methane	7500	U		15000	U
--2,4-Dichlorophenol	7500	U	IL	15000	U
--1,2,4-Trichlorobenzene	7500	U		15000	U
-Naphthalene	2500	J ✓	3.33	2700	DI 1L 1.5
-4-Chloroaniline	7500	U		15000	U
--Hexachlorobutadiene	7500	U		15000	U
-4-Chloro-3-Methylphenol	7500	U		15000	U
-2-Methylnaphthalene	790	J ✓		15000	U
--Hexachlorocyclopentadiene	7500	U		15000	U
--2,4,6-Trichlorophenol	7500	U		15000	U
-2,4,5-Trichlorophenol	18000	U		36000	U
-2-Chloronaphthalene	7500	U		15000	U
--2-Nitroaniline	18000	U		36000	U
-Dimethyl Phthalate	7500	U		15000	U
-Acenaphthylene	7500	U		15000	U
--2,6-Dinitrotoluene	7500	U		15000	U
--3-Nitroaniline	18000	U		36000	U
-Acenaphthene	9000	✓		10000	DI

FORM I SV-1

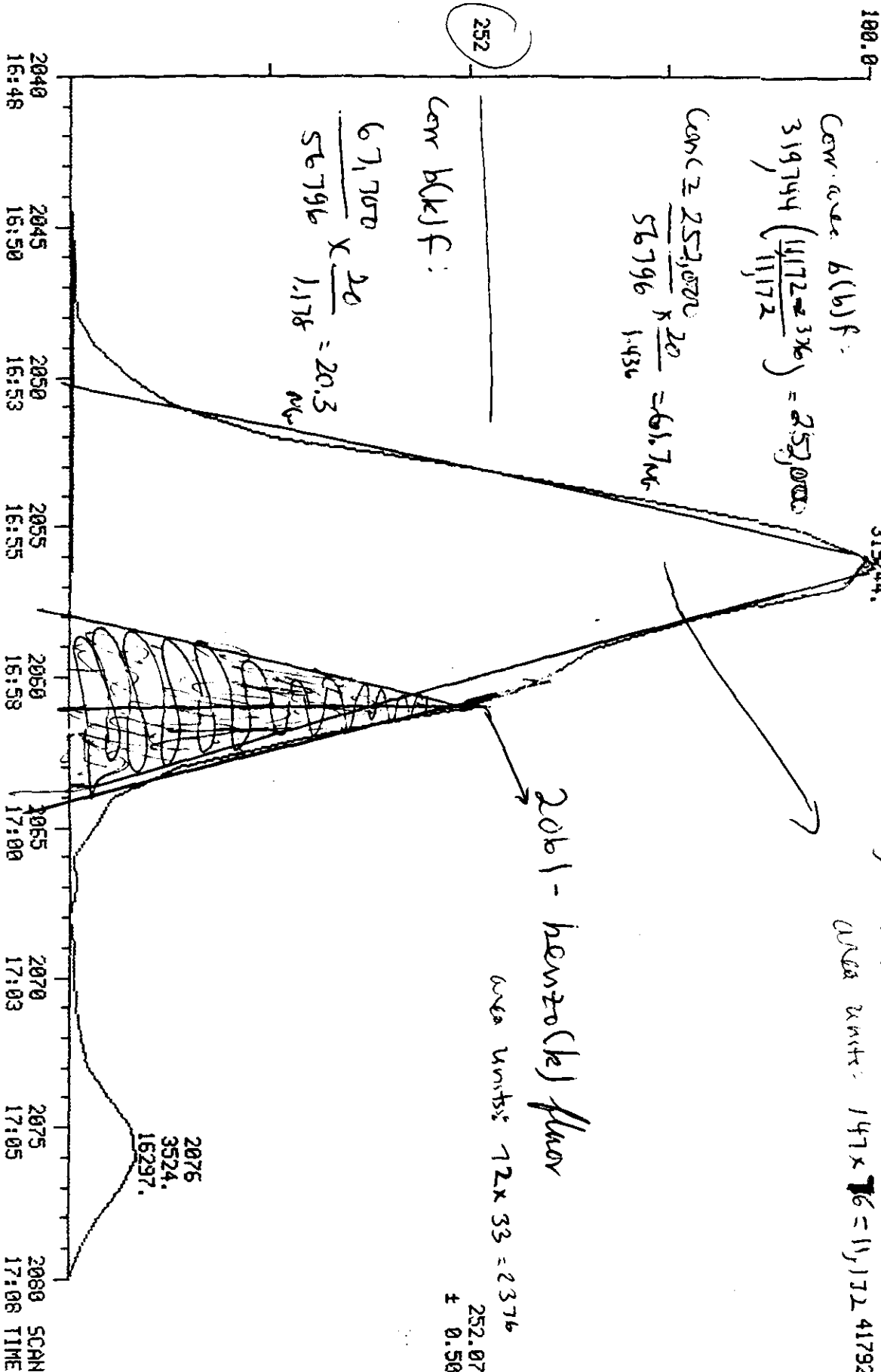
3/90

3/90

17744 CHY 11

1100

MIDMASS CHROMATOGRAM  
 02/11/92 3:44:00  
 SAMPLE: 2UL CC#477491 ID#CHY20 MS  
 COND.: EXTRACTED 2/4/92 1: 15 DILUTION  
 RANGE: G 1,2855 LABEL: N 1, 3.0 QUANT: A 1, 3.0 J 0 BASE: U 20, 3  
 DATA: GR077491C21 #1 SCANS 2040 TO 2080  
 CALI: GR077491C21 #2  
 CS#17744-CHY11 OMR21



SAMPLE DATA PACKAGE

17744 CHY 11

1148

56 \*459 D12-CHRYSENE (IS#5)  
 57 445 PYRENE (Q5#3) <129-00-0>  
 58 415 BUTYLBENZYL PHTHALATE (Q5#4) <85-68-7>  
 59 423 3,3'-DICHLOROBENZIDINE (Q5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (Q5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (Q5#7) <117-81-7>  
 62 418 CHRYSENE (Q5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (Q6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (Q6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (Q6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (Q6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (Q6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (Q6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (Q6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CHY 20

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1420 ✓	11:42	45	1.004	A BV	733607.	99.416 NG	11.72% <sup>YE</sup>
52	167	1455 ✓	11:59	45	1.028	A BV	53628.	14.904 NG	1.76% <sup>Y</sup>
53	178	1428 ✓	11:46	45	1.009	A VV	108429.	18.923 NG	2.23% <sup>Y</sup>
54	NOT FOUND								
55	202	1621	13:21	45	1.146	A BB	912634.	145.471 NG	17.15% <sup>YE</sup>
56	240	1864	15:21	56	1.000	A BV	155836.	20.000 NG	2.36% <sup>IS 5</sup>
57	202	1660	13:40	56	0.891	A BB	741494.	67.456 NG	7.95% <sup>Y</sup>
58	NOT FOUND								
59	NOT FOUND								
60	228	1863 ✓	15:21	56	0.999	A BV	372047.	43.356 NG	5.11% <sup>Y</sup>
61	149	1865	15:21	56	1.001	A BB	18959.	2.094 NG	0.25% <sup>Y</sup>
62	228	1869 ✓	15:23	56	1.003	A VV	274026.	42.252 NG	4.98% <sup>Y</sup>
63	264	2139	17:37	63	1.000	A BV	56796.	20.000 NG	2.36% <sup>IS 6</sup>
64	NOT FOUND								
65	252	2056	16:56	63	0.961	A BV	362987. <sup>252,000</sup>	81.647 NG <sup>46,000 ng/kg</sup>	9.62% <sup>Y</sup>

# SAMPLE DATA PACKAGE

17744 CHY 11

46,000 ng/kg 1116

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2061	16:56	63	0.961	A BV	362987. <sup>67,700</sup>	81.647 NG <sup>20.3 = 15,200 ng/kg</sup>	11.74% <sup>Y</sup>
67	252	2127	17:31	63	0.994	A VV	107344.	40.150 NG	4.73% <sup>Y</sup>
68	276	2451	20:11	63	1.146	A BB	37493.	14.028 NG	1.65% <sup>Y</sup>
69	278	2457	20:14	63	1.149	A BB	17340.	7.013 NG	0.83% <sup>Y</sup>
70	276	2544	20:57	63	1.189	A BB	25652.	9.986 NG	1.18% <sup>Y</sup>
71	112	499	4:07	1	0.732	A BV	11704.	3.325 NG	0.39% <sup>Y</sup>
72	99	632	5:12	1	0.927	A BV	9492.	2.260 NG	0.27% <sup>Y</sup>
73	132	650	5:21	1	0.953	A BB	7612.	2.227 NG	0.26% <sup>Y</sup>
74	152	705	5:48	1	1.034	A BB	4540.	1.818 NG	0.21% <sup>Y</sup>
75	82	767	6:19	13	0.873	A BV	6876.	1.487 NG	0.18% <sup>Y</sup>
76	172	1056	8:42	26	0.905	A BB	12948.	1.982 NG	0.23% <sup>Y</sup>
77	330	1299	10:42	26	1.113	A BB	2116.	1.366 NG	0.16% <sup>Y</sup>
78	244	1684	13:52	56	0.903	A BB	14052.	2.099 NG	0.25% <sup>Y</sup>

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:42	1.00	10.000	0.10	99.42	25.00	4.449	1.119	3.98
52	12:01	1.00	1.000	1.03	14.90	25.00	0.325	0.546	0.60
53	11:47	1.00	10.000	0.10	18.92	25.00	0.458	0.260	0.27

MIDMASS CHROMATOGRAM

02/13/92 1:22:00  
 SAMPLE: 2UL CC#477491 ID#CHY20 RE  
 COND.: EXTRACTED 2/4/92 1:30 DILUTION  
 RANGE: G 1, 2855 LABEL: N 1, 3.0 QUAN: A 1, 0.3 J 0 BASE: U 20, 3

*over*

DATA: GDJ77491A21 #1  
 CALL: GDJ77491A21 #2  
 CS#17744-CHY11  
 OMA21

SCANS 2015 TO 2125

19072. benzol(h) fluor

area units: 145 x 20 = 5776

$$\text{Conv. area: } 143716 \left( \frac{3710-652}{3776} \right) = 111,200$$

$$\text{Conv. conc. } \frac{111,200}{79552} \times \frac{20}{1.150} = 24.3 \text{ ng}$$

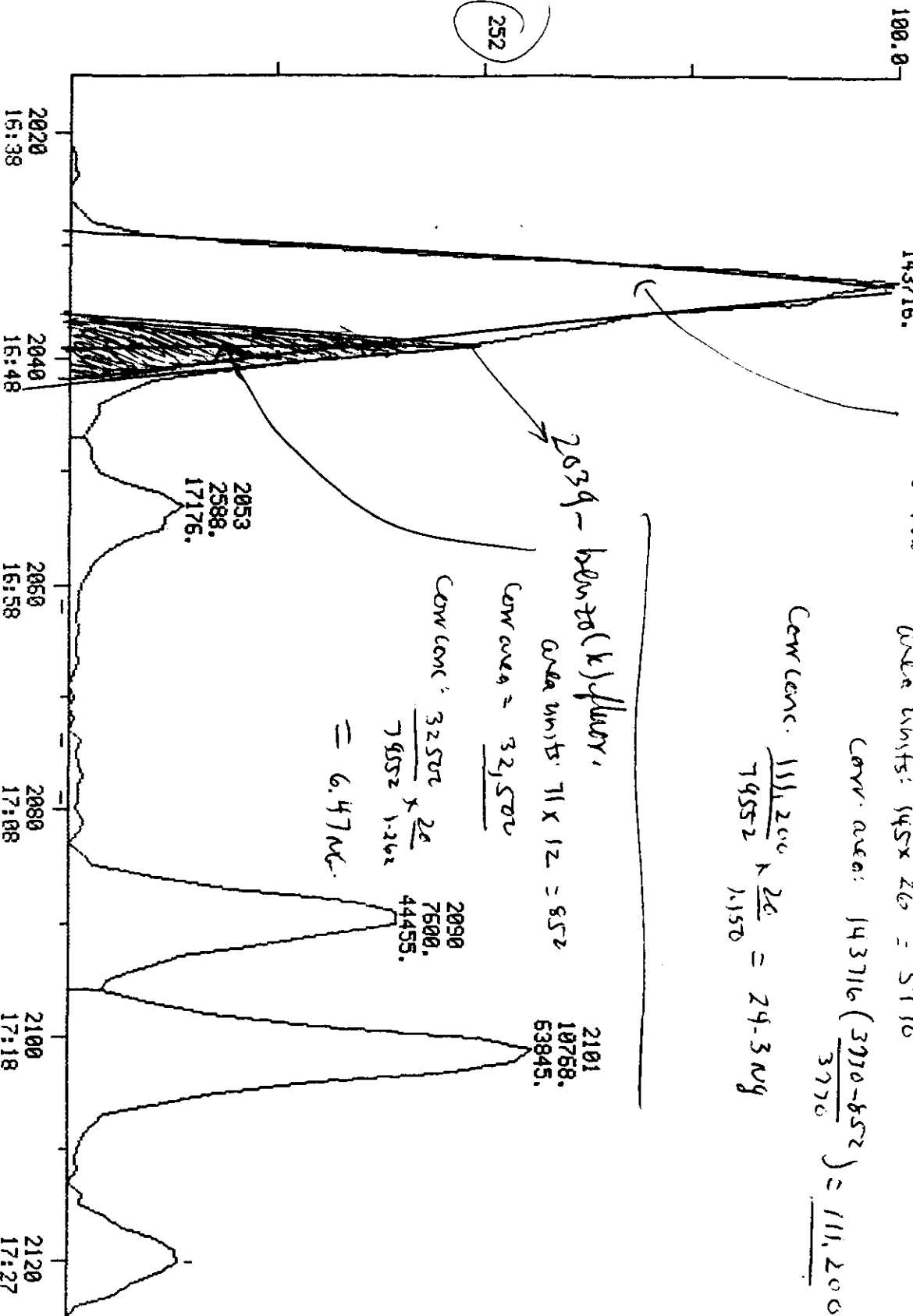
2039 - benzol(h) fluor

area units: 71 x 12 = 852

$$\begin{aligned} \text{Conv. area} &= 32,502 \\ \text{Conv. conc.} &= \frac{32,502}{79552} \times \frac{20}{1.262} \\ &= 6.47 \text{ ng} \end{aligned}$$

2090 7500. 44455.  
 2101 10768. 63845.

252.075  
 ± 0.500



60 405 BENZO(A)ANTHRACENE (Q5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (Q5#7) <117-81-7>  
 62 418 CHRYSENE (Q5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (Q6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (Q6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (Q6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (Q6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (Q6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (Q6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (Q6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CHY 20 2

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1405	11:34	45	1.003	A BV	469228.	34.065 NG	8.24 y D
52	167	1442	11:52	45	1.029	A BB	51260.	5.223 NG	1.26 y D
53	178	1413	11:38	45	1.009	A VB	106808.	8.912 NG	2.16 y D
54	NOT FOUND								
55	202	1605	13:13	45	1.146	A BB	477180.	47.539 NG	11.50 y D
56	240	1848	15:13	56	1.000	A BB	105528.	20.000 NG	4.84 y D
57	202	1643	13:32	56	0.889	A BB	370408.	35.704 NG	8.63 y D
58	149	1753	14:26	56	0.949	A BB	6680.	1.211 NG	0.29 y D
59	NOT FOUND								
60	228	1846	15:12	56	0.999	A BV	116968.	19.256 NG	4.66 y D
61	149	1851	15:15	56	1.002	A BB	14408.	1.719 NG	0.42 y D
62	228	1852	15:15	56	1.002	A VV	105604.	17.615 NG	4.26 y D
63	264	2113	17:24	63	1.000	A BB	79552.	20.000 NG	4.84 y D
64	NOT FOUND								
65	252	2033	16:45	63	0.962	A BV	142028.	24.315 NG	7.55 y D

111,200 24.315 NG = 36,5000/kg

# SAMPLE DATA PACKAGE

17744 CHY 11

1189

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2033	16:45	63	0.962	A BV	142028.	24.315 NG	6.88 y D
67	252	2101	17:18	63	0.994	A VV	66900.	16.362 NG	3.96 y D
68	276	2411	19:51	63	1.141	A BB	43812.	9.793 NG	2.37 y D
69	278	2417	19:54	63	1.144	A VB	14424.	3.967 NG	0.96 y D
70	276	2501	20:36	63	1.184	A BB	33040.	8.308 NG	2.01 y D
71	112	488	4:01	1	0.728	A BB	8272.	1.399 NG	0.34 y D
72	99	622	5:07	1	0.928	A BV	8324.	1.124 NG	0.27 y D
73	132	639	5:16	1	0.954	A BB	7112.	1.026 NG	0.25 y D
74	NOT FOUND								
75	82	757	6:14	13	0.873	A BB	9160.	1.189 NG	0.29 y D
76	172	1044	8:36	26	0.905	A BB	13160.	1.089 NG	0.26 y D
77	330	1286	10:35	26	1.115	A BB	2380.	1.142 NG	0.28 y D
78	244	1670	13:45	56	0.904	A BB	9896.	1.451 NG	0.35 y D

2039 6.47 x 1500 = 97000/kg

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:35	1.00	10.000	0.10	34.06	25.00	1.559	1.144	1.36
52	11:53	1.00	1.000	1.03	5.22	25.00	0.170	0.815	0.21
53	11:39	1.00	10.000	0.10	8.91	25.00	0.355	0.996	0.36
54	12:22		10.000						
55	13:14	1.00	10.000	0.11	47.54	25.00	1.586	0.834	1.90
56	15:14	1.00	10.000	0.10	20.00	20.00	1.000	1.000	1.00
57	13:32	1.00	10.000	0.08	35.70	25.00	1.428	1.266	1.12

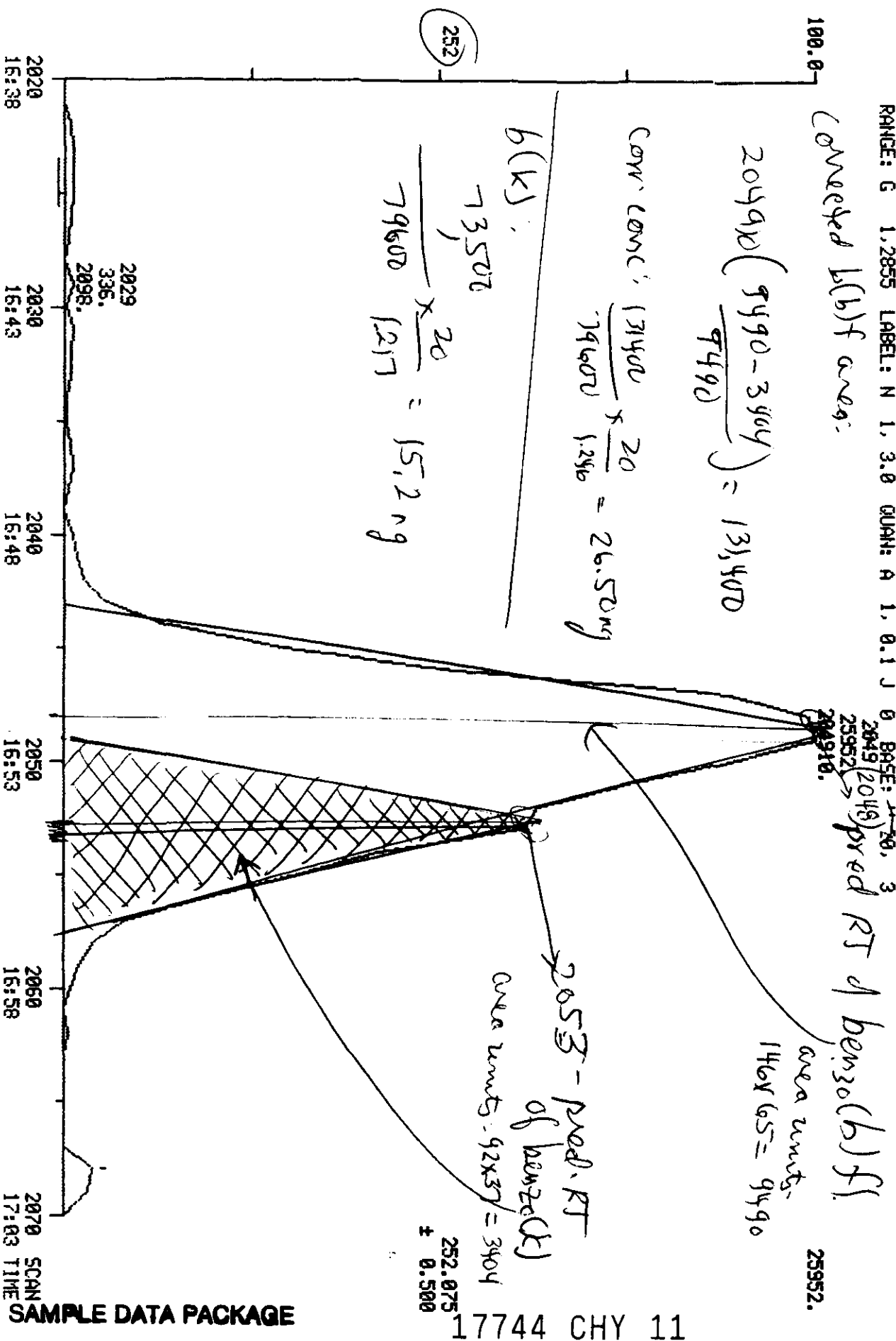


MIDMASS CHROMATOGRAM  
 02/12/92 4:43:08  
 SAMPLE: 2UL CC#477492 ID#CHY21 R5  
 CONDS.: EXTRACTED 2/4/92 1: 5 DILUTION  
 RANGE: G 1,2855 LABEL: N 1, 3.0 QUANT: A 1, 0.1 J 0  
 DATA: G2D77492A21 #1  
 CALL: G2D77492A21 #2  
 CS#17744-CHY11  
 OMA21  
 SCANS 2020 TO 2070

Converted b(b)f area:  
 20490 (9490 - 3404) = 13,400  
 14645 = 9490  
 area units  
 25952.  
 pred RT of benz(b)f

Cor. conc:  $\frac{13400}{19600} \times \frac{20}{1.246} = 26.50 \text{ ng}$

b(k):  
 $\frac{13500}{19600} \times \frac{20}{1.217} = 15.2 \text{ ng}$



65 407 BENZO(B)FLUORANTHENE (G6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (G6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (G6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (G6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (G6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (G6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #370 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

C4421

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1414	11:39	45	1.003	A VV	107240.	13.697 NG	3.4045
52	167	1451	11:57	45	1.029	A BB	12436.	2.816 NG	0.7045
53	178	1423	11:43	45	1.009	A VV	24560.	3.634 NG	0.9045
54	NOT FOUND								
55	202	1615	13:18	45	1.145	A BB	204028.	34.993 NG	8.6945
56	240	1858	15:18	56	1.000	A VB	115832.	20.000 NG	4.97 15.5
57	202	1653	13:37	56	0.890	A VB	164717.	13.861 NG	3.4445
58	149	1763	14:31	56	0.949	A BB	9848.	1.854 NG	0.4645
59	NOT FOUND								
60	228	1857	15:18	56	0.999	A BV	116709.	17.700 NG	4.4045
61	NOT FOUND								
62	228	1863	15:21	56	1.003	A VV	92921.	15.028 NG	3.7345
63	264	2132	17:33	63	1.000	A VV	79600.	20.000 NG	4.97 15.6
64	NOT FOUND								
65	252	2049	16:52	63	0.961	A BV	131400	26.50 NG	10.7345

# SAMPLE DATA PACKAGE

17744 CHY 11

1266

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2053	16:52	63	0.961	A BV	131400	26.50 NG	10.9945
67	252	2119	17:27	63	0.994	A VV	103751.	27.361 NG	6.7945
68	276	2438	20:05	63	1.144	A BB	59466.	15.128 NG	3.7645
69	278	2445	20:08	63	1.147	A*BV	21735.	6.763 NG	1.6845
70	NOT FOUND								
71	112	491	4:03	1	0.726	A BV	33444.	6.921 NG	1.72
72	99	626	5:09	1	0.926	A BV	37364.	6.348 NG	1.58
73	132	643	5:18	1	0.951	A BB	23572.	5.043 NG	1.25
74	152	699	5:45	1	1.034	A BB	14048.	4.080 NG	1.01
75	82	761	6:16	13	0.871	A BV	24972.	4.198 NG	1.04
76	172	1051	8:39	26	0.904	A BB	34916.	4.451 NG	1.11
77	330	1293	10:39	26	1.113	A BB	5480.	5.972 NG	1.48
78	244	1679	13:50	56	0.904	A VB	27628.	4.018 NG	1.00

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:41	1.00	10.000	0.10	13.70	25.00	0.626	1.142	0.55
52	11:59	1.00	1.000	1.03	2.82	25.00	0.073	0.644	0.11
53	11:45	1.00	10.000	0.10	3.63	25.00	0.143	0.986	0.15
54	12:28		10.000						
55	13:19	1.00	10.000	0.11	34.99	25.00	1.190	0.850	1.40
56	15:20	1.00	10.000	0.10	20.00	20.00	1.000	1.000	1.00
57	13:38	1.00	10.000	0.09	13.86	25.00	1.138	2.052	0.55
58	14:32	1.00	10.000	0.09	1.85	25.00	0.068	0.917	0.07
59	15:17		20.000						
60	15:19	1.00	10.000	0.10	17.70	25.00	0.806	1.138	0.71
61	15:21		10.000						
62	15:22	1.00	10.000	0.10	15.03	25.00	0.602	1.048	0.60

anal  
2/26/06

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

EPA SAMPLE

CHY22

CHY22RE

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11 : No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477493 : 477493

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GRJ77493A21 GRJ77493A05

Level: (low/med) LOW Date Received: 01/29/92 : 01/29/92

% Moisture: 20 decanted: (Y/N) N Date Extracted: 02/04/92 : 02/14/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92 : 02/24/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0 or: 1.0

GPC Cleanup: (Y/N) Y pH: 6.4

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

IS:  
3/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q	IS: <u>3/KG</u>	Q
108-95-2	Phenol	410	U	410	U
111-44-4	bis(2-Chloroethyl) Ether	410	U	410	U
95-57-8	2-Chlorophenol	410	U	410	U
541-73-1	1,3-Dichlorobenzene	410	U	410	U
106-46-7	1,4-Dichlorobenzene	410	U	410	U
95-50-1	1,2-Dichlorobenzene	410	U	410	U
95-48-7	2-Methylphenol	410	U	410	U
108-60-1	2,2'-Oxybis(1-Chloropropane)	410	U	410	U
106-44-5	4-Methylphenol	410	U	410	U
621-64-7	N-Nitroso-Di-n-Propylamine	410	U	410	U
67-72-1	Hexachloroethane	410	U	410	U
98-95-3	Nitrobenzene	410	U	410	U
78-59-1	Isophorone	410	U	410	U
88-75-5	2-Nitrophenol	410	U	410	U
105-67-9	2,4-Dimethylphenol	410	U	410	U
111-91-1	bis(2-Chloroethoxy) Methane	410	U	410	U
120-83-2	2,4-Dichlorophenol	410	U	410	U
120-82-1	1,2,4-Trichlorobenzene	410	U	410	U
91-20-3	Naphthalene	59	J	71	J ✓
106-47-8	4-Chloroaniline	410	U	410	U
87-68-3	Hexachlorobutadiene	410	U	410	U
59-50-7	4-Chloro-3-Methylphenol	410	U	410	U
91-57-6	2-Methylnaphthalene	75	J ✓	57	J
77-47-4	Hexachlorocyclopentadiene	410	U	410	U
88-06-2	2,4,6-Trichlorophenol	410	U	410	U
95-95-4	2,4,5-Trichlorophenol	980	U	990	U
91-58-7	2-Chloronaphthalene	410	U	410	U
88-74-4	2-Nitroaniline	980	U	990	U
131-11-3	Dimethyl Phthalate	410	U	410	U
208-96-8	Acenaphthylene	410	U	410	U
606-20-2	2,6-Dinitrotoluene	410	U	410	U
99-09-2	3-Nitroaniline	980	U	990	U
83-32-9	Acenaphthene	410	U	150	J ✓

FORM I SV-1

3/90

3/9

SAMPLE DATA PACKAGE

17744 CHY 11

1324

138

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

EPA SAMPLE

Lab Name: COMPUCHEM, RTP Contract: 68D10083

CHY22

CHY22RE

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

G No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477493

: 477493

Sample wt/vol: 30.5 (g/mL) G

Lab File ID: GRJ77493A21

GRJ77493AC

Level: (low/med) LOW

Date Received: 01/29/92

: 01/29/92

% Moisture: 20 decanted: (Y/N) N

Date Extracted: 02/04/92

d: 02/14/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 02/12/92

: 02/24/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

or: 1.0

GPC Cleanup: (Y/N) Y pH: 6.4

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG Q

TS:

G/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS:	Q	TS:	Q
		(ug/L or ug/Kg) <u>UG/KG</u>		<u>G/KG</u>	
51-28-5-----	2,4-Dinitrophenol	980	U	990	U
100-02-7-----	4-Nitrophenol	980	U	990	U
132-64-9-----	Dibenzofuran	410	U	61	J
121-14-2-----	2,4-Dinitrotoluene	410	U	410	U
84-66-2-----	Diethylphthalate	410	U	410	U
7005-72-3-----	4-Chlorophenyl-phenylether	410	U	410	U
86-73-7-----	Fluorene	410	U	89	J
100-01-6-----	4-Nitroaniline	980	U	990	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	980	U	990	U
86-30-6-----	N-Nitrosodiphenylamine (1)	410	U	410	U
101-55-3-----	4-Bromophenyl-phenylether	410	U	410	U
118-74-1-----	Hexachlorobenzene	410	U	410	U
87-86-5-----	Pentachlorophenol	980	U	990	U
85-01-8-----	Phenanthrene	370	J	900	J
120-12-7-----	Anthracene	55	J	150	J
86-74-8-----	Carbazole	68	J	120	J
84-74-2-----	Di-n-Butylphthalate	410	U	410	U
206-44-0-----	Fluoranthene	1100		1500	J
129-00-0-----	Pyrene	530		1300	J
85-68-7-----	Butylbenzylphthalate	410	U	410	U
91-94-1-----	3,3'-Dichlorobenzidine	410	U	410	U
56-55-3-----	Benzo(a)Anthracene	410		730	J
218-01-9-----	Chrysene	440		880	J
117-81-7-----	bis(2-Ethylhexyl)Phthalate	410	U	66	J
117-84-0-----	Di-n-Octyl Phthalate	410	U	410	U
205-99-2-----	Benzo(b)Fluoranthene	1200	X	1600	J
207-08-9-----	Benzo(k)Fluoranthene	<del>1300</del>	X	570	J
50-32-8-----	Benzo(a)Pyrene	380	J	550	
193-39-5-----	Indeno(1,2,3-cd)Pyrene	450		970	
53-70-3-----	Dibenz(a,h)Anthracene	98	J	180	J
191-24-2-----	Benzo(g,h,i)Perylene	350	J	810	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2 *Surrogate Monomers* 3/90

SAMPLE DATA PACKAGE

17744 CHY 11

*Similar*

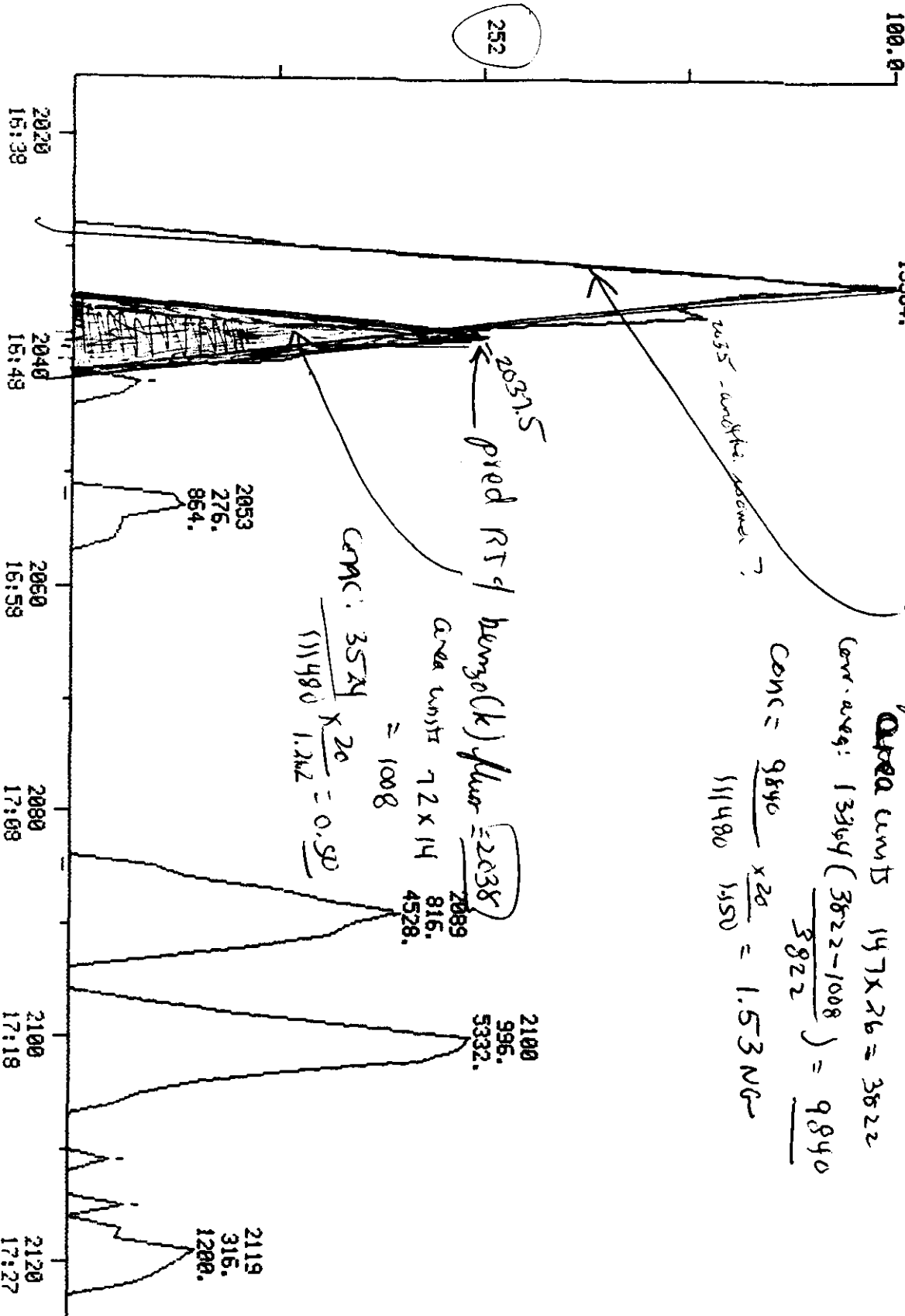
1325

3/

13

MIDMASS CHROMATOGRAM  
 02/12/92 22:47:00  
 SAMPLE: 2UL CC#477494 ID#CHY23 RE  
 CONDS.: EXTRACTED 2/4/92 UNDISTILLED  
 RANGE: G 1, 2855 LABEL: N 1, 3.0 QUANT: A 1, 0.3 J 0 BASE: U 20, 3

DATA: G2J77494A21 #1 SCANS 2015 TO 2125  
 CALL: G2J77494A21 #2  
 CS#17744-CHY11 QMA21



area units 147 x 26 = 3822  
 conc =  $\frac{9840}{11480} \times \frac{20}{1450} = 1.53 \text{ NG}$   
 conc =  $\frac{3524}{11480} \times \frac{20}{1.262} = 0.50$

59 423 3,3'-DICHLOROBENZIDINE (G5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (G5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (G5#7) <117-81-7>  
 62 418 CHRYSENE (G5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (G6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (G6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (G6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (G6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (G6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (G6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (G6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CHY 23

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	NOT FOUND	1804					1804	0.579	0.31
52	NOT FOUND							0.202	
53	178	1404	11:34	45	1.003	A BB	18052	1.032 NG	0.31
54	NOT FOUND								
55	202	1604	13:13	45	1.146	A BB	27708	1.891 NG	0.57
56	240	1847	15:13	56	1.000	A BV	152580	20.000 NG	5.99
57	202	1642	13:31	56	0.889	A BB	22528	1.502 NG	0.45
58	NOT FOUND								
59	NOT FOUND								
60	228	1851	15:15	56	1.002	A VB	9116	1.038 NG	0.31
61	NOT FOUND								
62	228	1851	15:15	56	1.002	A VB	9116	1.052 NG	0.32
63	264	2113	17:24	63	1.000	A BB	111480	20.000 NG	5.99
64	NOT FOUND								
65	252	2033	16:45	63	0.962	A BB	9840	1.53 x 4 = 64 ug/kg	0.62

# SAMPLE DATA PACKAGE

17744 CHY 11

1471

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2033	16:45	63	0.962	A BB	3524	0.50	0.57
67	NOT FOUND								
68	NOT FOUND								
69	NOT FOUND								
70	NOT FOUND								
71	112	487	4:01	1	0.727	A BB	248420	29.167 NG	8.74
72	99	620	5:06	1	0.925	A BV	298432	27.974 NG	8.38
73	132	637	5:15	1	0.951	A BB	230712	23.107 NG	6.92
74	152	692	5:42	1	1.033	A BB	141232	21.316 NG	6.39
75	82	754	6:13	13	0.871	A BV	256564	23.554 NG	7.06
76	172	1043	8:35	26	0.905	A BB	377756	22.372 NG	6.70
77	330	1284	10:34	26	1.114	A BB	74604	25.642 NG	7.68
78	244	1669	13:45	56	0.904	A BV	243412	24.690 NG	7.40

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:35		10.000						
52	11:53		1.000						
53	11:39	0.99	10.000	0.10	1.03	25.00	0.041	0.996	0.04
54	12:22		10.000						
55	13:14	1.00	10.000	0.11	1.89	25.00	0.063	0.834	0.08
56	15:14	1.00	10.000	0.10	20.00	20.00	1.000	1.000	1.00
57	13:32	1.00	10.000	0.08	1.50	25.00	0.050	0.750	0.05

59 423 3,3'-DICHLOROBENZIDINE (Q5#5) <91-94-1>  
 60 405 BENZO(A)ANTHRACENE (Q5#6) <56-55-3>  
 61 413 BIS(2-ETHYLHEXYL) PHTHALATE (Q5#7) <117-81-7>  
 62 418 CHRYSENE (Q5#8) <218-01-9>  
 63 \*497 D12-PERYLENE (IS#6)  
 64 429 DI-N-OCTYL PHTHALATE (Q6#2) <117-84-0>  
 65 407 BENZO(B)FLUORANTHENE (Q6#3) <205-99-2>  
 66 409 BENZO(K)FLUORANTHENE (Q6#4) <207-08-9>  
 67 406 BENZO(A)PYRENE (Q6#5) <50-32-8>  
 68 437 INDENO(1,2,3-C,D)PYRENE (Q6#6) <193-39-5>  
 69 419 DIBENZO(A,H)ANTHRACENE (Q6#7) <53-70-3>  
 70 408 BENZO(G,H,I)PERYLENE (Q6#8) <191-24-2>  
 71 #619 2-FLUOROPHENOL (SS#1)  
 72 #612 D5-PHENOL (SS#2)  
 73 #634 2-CHLOROPHENOL-D4 (SS#3)  
 74 #570 1,2-DICHLOROBENZENE-D4 (SS#4)  
 75 #447 D5-NITROBENZENE (SS#5)  
 76 #448 2-FLUOROBIPHENYL (SS#6)  
 77 #628 2,4,6-TRIBROMOPHENOL (SS#7)  
 78 #496 D14-TERPHENYL (SS#8)

CH426

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
51	178	1413	11:38	45	1.003	A BV	15848.	1.594 NG	0.34 yes
52	NOT FOUND								
53	178	1413	11:38	45	1.003	A BV	15848.	1.847 NG	0.39 NO
54	NOT FOUND								
55	202	1613	13:17	45	1.145	A BB	26428.	3.570 NG	0.76 yes
56	240	1857	15:18	56	1.000	A BB	106520.	20.000 NG	4.24 5.5
57	202	1652	13:36	56	0.890	A BB	22416.	2.051 NG	0.43 yes
58	NOT FOUND								
59	NOT FOUND								
60	228	1861	15:20	56	1.002	A VB	10132.	1.782 NG	0.35 yes
61	NOT FOUND								
62	228	1861	15:20	56	1.002	A VB	10132.	1.782 NG	0.38 yes
63	264	2130	17:32	63	1.000	A BB	82768.	20.000 NG	4.24 ES6
64	NOT FOUND								
65	252	2046	16:51	63	0.961	A BB	17892.	3.467 NG	0.74 yes

# SAMPLE DATA PACKAGE

17744 CHY 11

2050

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hght)	Amount	%Tot
66	252	2046	16:51	63	0.961	A BB	17892.	3.467 NG	0.75 yes
67	252	2117	17:26	63	0.994	A VB	7012.	1.778 NG	0.38 yes
68	276	2437	20:04	63	1.144	A BB	5324.	1.303 NG	0.28 NO
69	NOT FOUND								
70	NOT FOUND								
71	112	491	4:03	1	0.728	A BV	235688.	46.665 NG	9.89
72	99	624	5:08	1	0.926	A BV	312132.	50.732 NG	10.76
73	132	642	5:17	1	0.953	A BV	184024.	37.666 NG	7.99
74	152	697	5:44	1	1.034	A BB	111300.	30.926 NG	6.56
75	82	759	6:15	13	0.870	A BV	229236.	35.780 NG	7.59
76	172	1050	8:39	26	0.904	A BB	293684.	32.373 NG	6.86
77	330	1292	10:38	26	1.113	A BB	44132.	41.592 NG	8.82
78	244	1678	13:49	56	0.904	A BB	222416.	35.176 NG	7.46

No	Ret(L)	Ratio	RRT(L)	Ratio	Amnt	Amnt(L)	R. Fac	R. Fac(L)	Ratio
51	11:41	1.00	10.000	0.10	1.59	25.00	0.073	1.142	0.06
52	11:59		1.000						
53	11:45	0.99	10.000	0.10	1.85	25.00	0.073	0.986	0.07
54	12:28		10.000						
55	13:19	1.00	10.000	0.11	3.57	25.00	0.121	0.850	0.14

7C  
SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
 Instrument ID: OWA08 Calibration date: 02/17/92 Time: 1746  
 Lab File ID: HG920217B08 Init. Calib. Date(s): 01/07/92 01/07/92  
 Init. Calib. Times: 1446 1920

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Diethylphthalate	1.602	1.594		0.5	
4-Chlorophenyl-phenylether	0.618	0.718	0.400	-16.2	25.0
Fluorene	1.181	1.249	0.900	-5.8	25.0
4-Nitroaniline	0.224	0.251		-12.0	
4,6-Dinitro-2-Methylphenol	0.121	0.134		-10.7	
N-Nitrosodiphenylamine (1)	0.464	0.423		8.8	
4-Bromophenyl-phenylether	0.242	0.262	0.100	-8.3	25.0
Hexachlorobenzene	0.336	0.366	0.100	-8.9	25.0
Pentachlorophenol	0.162	0.163	0.050	-0.6	25.0
Phenanthrene	1.077	1.079	0.700	-0.2	25.0
Anthracene	0.971	0.967	0.700	0.4	25.0
Carbazole	0.753	0.897		-19.1	
Di-n-Butylphthalate	1.782	1.674		6.1	
Fluoranthene	0.870	1.041	0.600	-19.7	25.0
Pyrene	1.484	1.339	0.600	9.8	25.0
Butylbenzylphthalate	1.103	0.863		21.8	
3,3'-Dichlorobenzidine	0.202	0.191		5.4	
Benzo(a)Anthracene	1.122	1.163	0.800	-3.7	25.0
Chrysene	1.047	1.073	0.700	-2.5	25.0
bis(2-Ethylhexyl) Phthalate	1.793	1.358		24.3	
Di-n-Octyl Phthalate	3.217	2.890		10.2	
Benzo(b)Fluoranthene	1.268	1.341	0.700	-5.8	25.0
Benzo(k)Fluoranthene	1.164	1.150	0.700	1.2	25.0
Benzo(a)Pyrene	1.034	1.049	0.700	-1.5	25.0
Indeno(1,2,3-cd)Pyrene	1.142	0.897	0.500	21.4	25.0
Dibenz(a,h)Anthracene	0.961	0.781	0.400	18.7	25.0
Benzo(g,h,i)Perylene	1.033	0.761	0.500	26.3	25.0
Nitrobenzene-d5	0.486	0.516	0.200	-6.2	25.0
2-Fluorobiphenyl	1.308	1.354	0.700	-3.5	25.0
Terphenyl-d14	1.104	1.033	0.500	6.4	25.0
Phenol-d5	1.855	2.007	0.800	-8.2	25.0
2-Fluorophenol	1.513	1.600	0.600	-5.8	25.0
2,4,6-Tribromophenol	0.224	0.266		-18.8	
2-Chlorophenol-d4	1.396	1.536	0.800	-10.0	25.0
1,2-Dichlorobenzene-d4	0.847	0.908	0.400	-7.2	25.0

(1) Cannot be separated from Diphenylamine  
 All other compounds must meet a minimum RRF of 0.010.

*affects Chy 16*



ORIGINAL

iPod

2E  
WATER PESTICIDE SURROGATE RECOVERY

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02GC Column (1): DB-608 ID: 0.53(mm) GC Column (2): RTX-1701 ID: 0.53(mm)

	EPA SAMPLE NO.	TCX 1 %REC #	TCX 2 %REC #	DCB 1 %REC #	DCB 2 %REC #	OTHER (1)	OTHER (2)	TOT OUT
01	CHY07MS	89	94	100	110			0
02	CHY07MSD	91	93	110	110			0
03	CHY02	84	86	100	110			0
04	CHY03	89	91	100	110			0
05	CHY04	88	89	100	110			0
06	CHY05	82	77	100	110			0
07	CHY06	80	74	100	100			0
08	CHY07	74	83	95	100			0
09	CHY08	79	78	97	100			0
10	CHY09	82	79	99	100			0
11	CHY10	80	77	100	110			0
12	CHY24	79	95	100	110			0
13	PBLK08	79	82	91	97			0
14								
15								
16								
17								
18								
19								
20								
21								
22								
23								
24								
25								
26								
27								
28								
29								
30								

QC LIMITS

TCX = Tetrachloro-m-xylene (60-150)  
DCB = Decachlorobiphenyl (60-150)

# Column to be used to flag recovery values.  
\* Values outside of QC Limits  
D Surrogate diluted out.

3E  
WATER PESTICIDE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6759HQ SDG No.: CHY02

Matrix Spike - EPA Sample No.: CHY07

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)	MS CONCENTRATION (ug/L)	MS % REC #	QC. LIMITS REC.
=====	=====	=====	=====	=====	=====
gamma-BHC (Lindane)	0.500	0.00	0.42	84	56-123
Heptachlor	0.500	0.00	0.41	82	40-131
Aldrin	0.500	0.00	0.39	78	40-120
Dieldrin	1.000	0.00	0.83	83	52-126
Endrin	1.000	0.00	0.93	93	56-121
4,4'-DDT	1.000	0.00	0.90	90	38-127

COMPOUND	SPIKE ADDED (ug/L)	MSD CONCENTRATION (ug/L)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
=====	=====	=====	=====	=====	=====	=====
gamma-BHC (Lindane)	0.500	0.42	84	0	15	56-123
Heptachlor	0.500	0.41	82	0	20	40-131
Aldrin	0.500	0.39	78	0	22	40-120
Dieldrin	1.000	0.84	84	1	18	52-126
Endrin	1.000	0.95	95	2	21	56-121
4,4'-DDT	1.000	0.92	92	2	27	38-127

# Column to be used to flag recovery and RPD values with an asterisk

\* Values outside of QC limits

RPD: 0 out of 6 outside limits

Spike Recovery: 0 out of 12 outside limits

COMMENTS: \_\_\_\_\_

2F  
SOIL PESTICIDE SURROGATE RECOVERY

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

GC Column (1): DB-608 ID: 0.53(mm) GC Column (2): RTX-1701 ID: 0.53(mm)

	EPA SAMPLE NO.	TCX 1 %REC #	TCX 2 %REC #	DCB 1 %REC #	DCB 2 %REC #	OTHER (1)	OTHER (2)	TOT OUT
01	CHY11	140	0*	63	74			1
02	CHY11DL	140	48*	48*	65			2
03	CHY12	52*	55*	56*	74			3
04	CHY13	80	120	110	99			0
05	CHY14	100	0*	46*	64			2
06	CHY14DL	99	0*	70	57*			2
07	CHY16	95	85	340*	380*			2
08	CHY17	76	60	0*	340*			2
09	CHY19	96	120	100	130			0
10	CHY20	65	41*	280*	280*			3
11	CHY22 *	25*	81	43*	140			2
12	CHY23	77	97	57*	150			1
13	CHY26	78	88	63	110			0
14	PBLK18	85	100	97	100			0
15	CHY18MS	68	52*	200*	220*			3
16	CHY18MSD	74	91	310*	150			1
17	CHY18	84	110	180*	190*			2
18	PBLK62	75	94	74	77			0
19	CHY15	0D	0D	0D	0D			0
20	CHY15DL	0D	0D	0D	0D			0
21	CHY21	53*	37*	300*	270*			4
22	PBLK73	74	89	83	81			0
23								
24								
25								
26								
27								
28								
29								
30								

\* CHY22 - no obvious matrix effects to account for variable TCX recoveries

QC LIMITS  
TCX = Tetrachloro-m-xylene (60-150)  
DCB = Decachlorobiphenyl (60-150)

# Column to be used to flag recovery values.

\* Values outside of QC Limits

D Surrogate diluted out.

3F  
SOIL PESTICIDE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix Spike - EPA Sample No.: CHY18

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC. LIMITS REC.
=====	=====	=====	=====	=====	=====
gamma-BHC (Lindane)	23.148	0.00	15	65	46-127
Heptachlor	23.148	0.00	15	65	35-130
Aldrin	23.148	0.36	14	59	34-132
Dieldrin	46.296	0.27	29	62	31-134
Endrin	46.296	2.6	36	72	42-139
4,4'-DDT	46.296	0.48	29	62	23-134

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
=====	=====	=====	=====	=====	=====	=====
gamma-BHC (Lindane)	23.148	15	65	0	50	46-127
Heptachlor	23.148	14	60	8	31	35-130
Aldrin	23.148	14	59	0	43	34-132
Dieldrin	46.296	29	62	0	38	31-134
Endrin	46.296	34	68	6	45	42-139
4,4'-DDT	46.296	32	68	9	50	23-134

# Column to be used to flag recovery and RPD values with an asterisk

\* Values outside of QC limits

RPD: 0 out of 6 outside limits

Spike Recovery: 0 out of 12 outside limits

COMMENTS: \_\_\_\_\_

ORIGINAL  
(FBI)

LABORATORY NOTICE  
Pesticide Fraction

Sample I.D.'s 477485 / 477492 Client I.D.'s CHY15 / CHY21  
Case 17744 SDG CHY11 SAS 65794Q

The above samples were extracted and analyzed within contractual holding times. After all the samples in the extraction batch had been completed, the GC/MS data reviewers sent a TIC report to the laboratory requesting confirmation of PCB in sample I.D. #477485 (CHY15), which did not match the I.D. number of the sample sent for confirmation prior to their report, which was sample I.D. #477492 (CHY21). These samples were then reextracted outside holding times and reanalyzed. The reextract data matched the TIC report for CHY15, however, we could not verify the original results for all of the target analytes in the reextract data for CHY21. We are, therefore, reporting the reextract data for these two samples with reference to this notice.

Pesticide fraction of  
Sample CHY15 and  
CHY21 were  
initially switched.  
Rg 6/7/92

Data Reviewer/I.D. (b) (4) / 1569  
Date 2/26/92



COMPUCHEM  
LABORATORIES, INC.

P.O. Box 12652 3308 Chapel Hill/Nelson Highway Research Triangle Park, NC 27709 (919) 549-8263

SDG Narrative # CHY11  
Case # 17744 SAS # 6579HQ  
Contract # SAS6579HQ  
CompuChem Laboratories, Inc.

SAMPLE IDENTIFICATIONS: CHY11, CHY12, CHY13, CHY14, CHY15,  
CHY16, CHY17, CHY18, CHY19, CHY20, CHY21, CHY22, CHY23, CHY26

All pertinent Quality Assurance and Laboratory notices for CASE# 17744, SDG# CHY11, SAS# 6579HQ are included in the sample data package.

**PESTICIDES:**

Pesticides Target Compound List (TCL) analytes were confirmed in nine of these samples above the Contract Required Quantitation Limits (CRQLs). Holding time requirements were initially met for all of these samples.

The initial analyses of samples CHY11 and CHY14 yielded TCL analytes that were above the analytical range. Each of these samples were reanalyzed at a 2X dilution in order to report the analytes within the analytical range. Both analyses for each of these samples are reported as billable.

Semivolatile analysis of sample CHY15 indicated that a PCB confirmational analysis was necessary. Inspection of the pesticide data for this sample indicated that no PCBs were present in this sample; however, the pesticide analysis of sample CHY21 contained PCB. These samples were reextracted outside of holding time and reanalyzed. Results from the reanalysis of sample CHY15 verified the semivolatile GCMS TIC identification of PCBs in the sample. The original TCL results of CHY21 could not be verified. The reextract data which was outside of holding time is being reported with this case.

The surrogate standard recovery of decachlorobiphenyl (DCB) was outside of the control range on both columns in the analysis of samples CHY18MS, CHY16, CHY17, CHY18, CHY20, and CHY21. These outliers have been attributed to matrix interference and not to deficiencies in laboratory practice.

The surrogate standard recovery of DCB was outside of the control range on the DB-608 column in the analysis of sample CHY18MSD. The surrogate standard recovery of tetrachloro-m-xylene was outside of the control range on the RTX-1701 column in the analyses of samples CHY11, CHY14, and CHY14DL. These outliers are due to matrix interferences which affected accurate quantitation.

No surrogate recoveries are reported in sample CHY15 and CHY15DL due to the 20X and 200X dilution that were necessary to report accurate results. Arochlor 1260 was detected in the GC analysis at concentrations that require GCMS analysis. Results from the GCMS qualitative analysis positively identified arochlor 1260 in sample CHY15. Both the 20X and the 200X dilution are reported as billable.

*Handwritten:* Pesticide fraction of samples initially switched  
pg 6/7/92

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY21 RE

Lab Name: COMPUCHEM, RTPContract: 68D10083Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11Matrix: (soil/water) SOILLab Sample ID: 477492 R1Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 35 decanted: (Y/N) NDate Received: 01/29/92Extraction: (SepF/Cont/Sonc) SONCDate Extracted: 02/20/92Concentrated Extract Volume: 5000(uL)Date Analyzed: 02/22/92Injection Volume: 2.0(uL)Dilution Factor: 1GPC Cleanup: (Y/N) XpH: 6.6Sulfur Cleanup: (Y/N) N

Reanal. → no PCBs at High Level Q

CAS NO.

COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

319-84-6-----	alpha-BHC	2.6	U
319-85-7-----	beta-BHC	2.6	U
319-86-8-----	delta-BHC	2.6	U
58-89-9-----	gamma-BHC (Lindane)	0.18	JP
76-44-8-----	Heptachlor	0.81	JPB
309-00-2-----	Aldrin	1.4	JP
1024-57-3-----	Heptachlor epoxide	2.6	U
959-98-8-----	Endosulfan I	2.6	U
60-57-1-----	Dieldrin	2.4	JP
72-55-9-----	4,4'-DDE	6.4	
72-20-8-----	Endrin	5.1	U
33213-65-9-----	Endosulfan II	5.1	U
72-54-8-----	4,4'-DDD	5.1	U
1031-07-8-----	Endosulfan sulfate	5.1	U
50-29-3-----	4,4'-DDT	13	P
72-43-5-----	Methoxychlor	110	P
53494-70-5-----	Endrin ketone	5.1	U
7421-93-4-----	Endrin aldehyde	20	P
5103-71-9-----	alpha-Chlordane	2.4	JP
5103-74-2-----	gamma-Chlordane	5.7	P
8001-35-2-----	Toxaphene	260	U
12674-11-2-----	Aroclor-1016	51	U
11104-28-2-----	Aroclor-1221	100	U
11141-16-5-----	Aroclor-1232	51	U
53469-21-9-----	Aroclor-1242	51	U
12672-29-6-----	Aroclor-1248	51	U
11097-69-1-----	Aroclor-1254	51	U
11096-82-5-----	Aroclor-1260	460	PB

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CHY21 *initial*

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477492 D51

Sample wt/vol: 30.10(g/ml)G

Lab File ID: PC077492B20

% Moisture: 35 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 50

GPC Cleanup: (Y/N) Y

pH: 6.6

Sulfur Cleanup: (Y/N) N

*initial - high level PCBs*

CAS NO.

COMPOUND

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

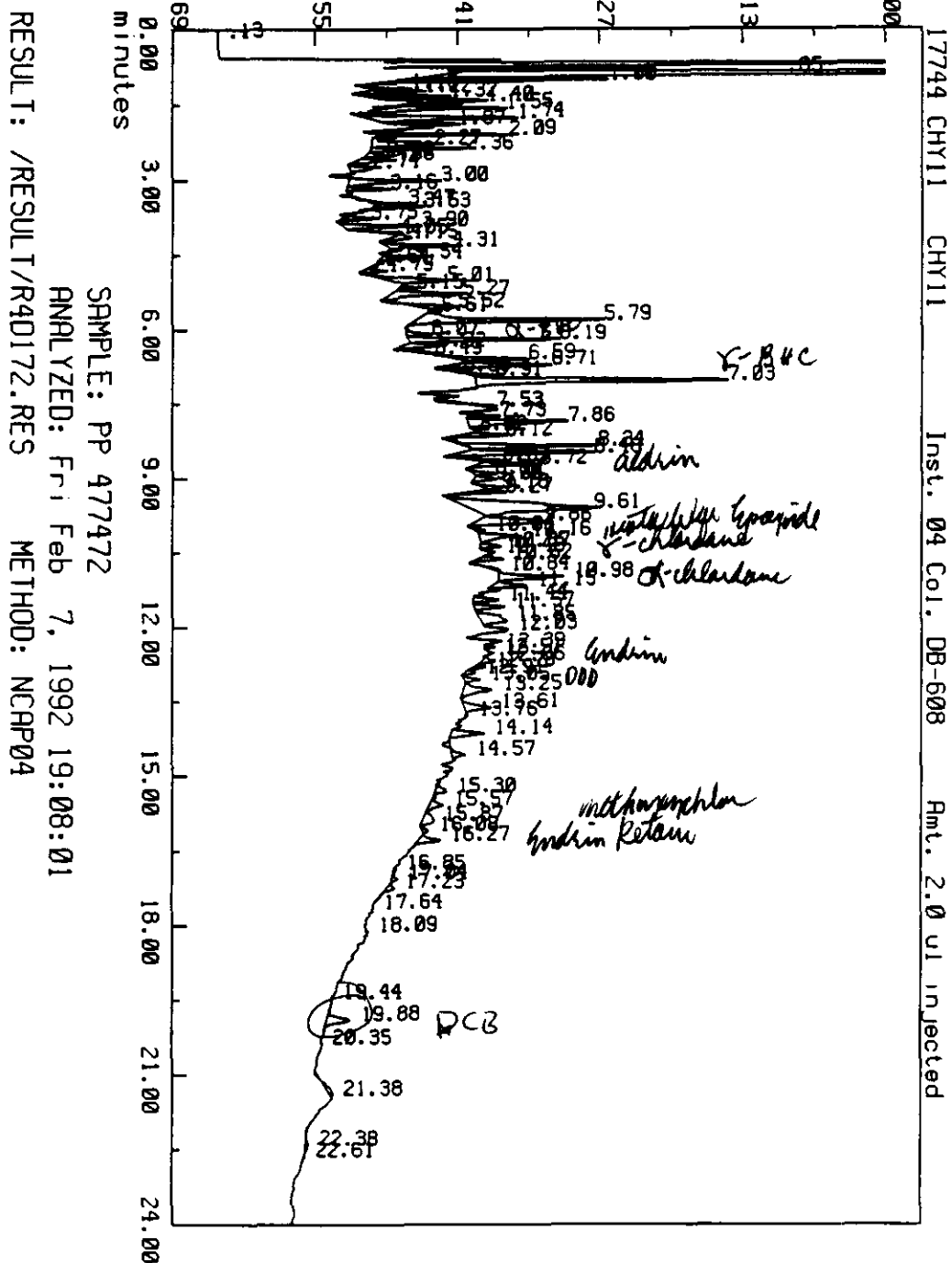
Q

319-84-6-----	alpha-BHC	48	JP
319-85-7-----	beta-BHC	130	U
319-86-8-----	delta-BHC	130	U
58-89-9-----	gamma-BHC (Lindane)	130	U
76-44-8-----	Heptachlor	130	U
309-00-2-----	Aldrin	94	JP
1024-57-3-----	Heptachlor epoxide	130	U
959-98-8-----	Endosulfan I	130	U
60-57-1-----	Dieldrin	250	U
72-55-9-----	4,4'-DDE	250	U
72-20-8-----	Endrin	250	U
33213-65-9-----	Endosulfan II	250	U
72-54-8-----	4,4'-DDD	250	U
1031-07-8-----	Endosulfan sulfate	250	U
50-29-3-----	4,4'-DDT	250	U
72-43-5-----	Methoxychlor	1300	U
53494-70-5-----	Endrin ketone	250	U
7421-93-4-----	Endrin aldehyde	250	U
5103-71-9-----	alpha-Chlordane	130	U
5103-74-2-----	gamma-Chlordane	130	U
8001-35-2-----	Toxaphene	13000	U
12674-11-2-----	Aroclor-1016	2500	U
11104-28-2-----	Aroclor-1221	5100	U
11141-16-5-----	Aroclor-1232	2500	U
53469-21-9-----	Aroclor-1242	2500	U
12672-29-6-----	Aroclor-1248	2500	U
11097-69-1-----	Aroclor-1254	2500	U
11096-82-5-----	Aroclor-1260	660000	C

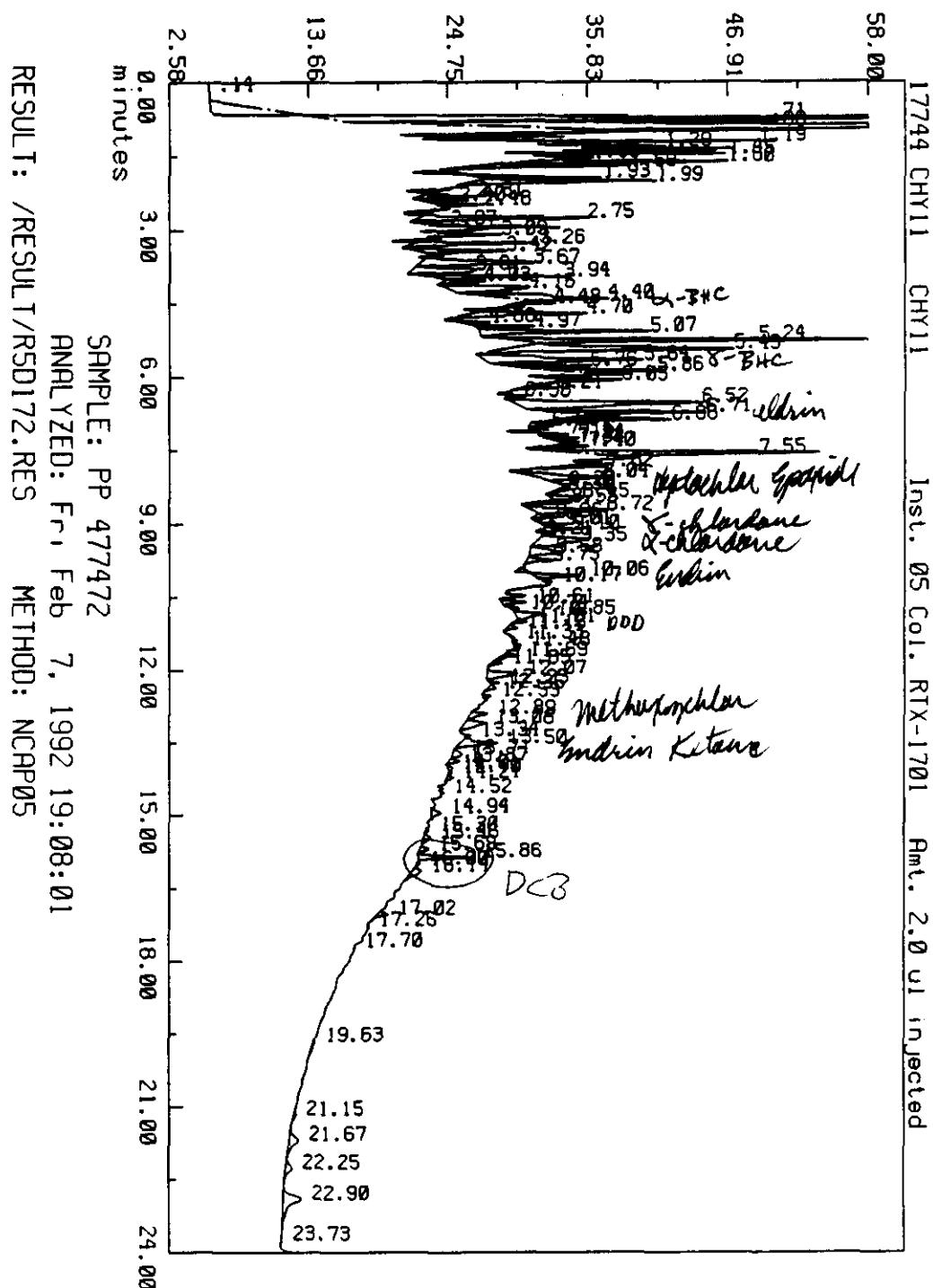


Example

Chromatogram illustrating closely-spaced peaks with continuous low-level interferences at same concentration range as targets.



AMPLITUDE/1000  
Force Normalized  
( 5.50, 58.00)



10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY11DL

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477472 D20

Date(s) Analyzed: 02/13/92 02/13/92

Instrument ID (1): VARIAN05

Instrument ID (2): VARIAN04

GC Column(1): RTX-1701

ID: 0.53(mm)

GC Column(2): DB-608

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM TO	CONCENTRATION	%D
gamma-BHC (Lindane)	1	5.65	5.61 5.71	31	
	2	7.03	6.98 7.08	67	116.1
delta-BHC	1	7.75	7.73 7.83	1.5	
	2	8.12	8.12 8.22	12	700.0
Aldrin	1	6.71	6.63 6.73	14	
	2	8.63	8.60 8.70	3.7	278.4
Heptachlor epoxide	1	8.20	8.16 8.30	1.7	
	2	10.15	10.05 10.19	9.5	458.8
gamma-Chlordane	1	9.01	8.87 9.01	3.3	
	2	10.48	10.47 10.61	2.7	22.2
alpha-Chlordane	1	9.10	9.06 9.20	4.4	
	2	10.97	10.88 11.02	12	172.7
Endrin	1	10.06	10.03 10.17	6.7	
	2	12.66	12.63 12.77	3.9	71.8
4,4'-DDD	1	11.13	11.07 11.21	1.8	
	2	13.04	13.02 13.16	4.8	166.7

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY11

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477472

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608 ID: 0.53(mm) GC Column(2): RTX-1701 ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
alpha-BHC	1	6.07	6.00	6.10	3.7	
	2	4.88	4.79	4.89	3.4	8.8
gamma-BHC (Lindane)	1	7.03	7.00	7.10	52	
	2	5.64	5.63	5.73	34	52.9
Aldrin	1	8.63	8.62	8.72	4.4	
	2	6.71	6.64	6.74	17	286.4
Heptachlor epoxide	1	10.16	10.06	10.20	11	
	2	8.20	8.17	8.31	0.95	999.9
gamma-Chlordane	1	10.48	10.48	10.62	1.0	
	2	9.01	8.89	9.03	4.1	310.0
alpha-Chlordane	1	10.98	10.90	11.04	14	
	2	9.10	9.07	9.21	5.3	164.2
Endrin	1	12.66	12.66	12.80	3.2	
	2	10.06	10.06	10.20	9.8	206.2
4,4'-DDD	1	13.05	13.03	13.17	5.4	
	2	11.13	11.08	11.22	2.8	92.9

page 1 of 2

SAMPLE DATA PACKAGE

FORM X PEST-1

17744 CHY 11

3/90

2251

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY12

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477476

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608

ID: 0.53(mm)

GC Column(2): RTX-1701

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
alpha-BHC	1	6.07	6.00	6.10	0.38	
	2	4.87	4.79	4.89	0.48	26.3
Heptachlor	1	7.86	7.80	7.90	1.8	
	2	6.08	6.08	6.18	0.13	999.9
Aldrin	1	8.65	8.62	8.72	1.1	
	2	6.72	6.64	6.74	2.3	109.1
gamma-Chlordane	1	10.55	10.48	10.62	4.5	
	2	8.94	8.89	9.03	3.5	28.6
alpha-Chlordane	1	10.93	10.90	11.04	4.1	
	2	9.13	9.07	9.21	9.3	126.8
4,4'-DDE	1	11.66	11.61	11.75	4.4	
	2	9.36	9.31	9.45	9.4	113.6
Endrin	1	12.70	12.66	12.80	7.6	
	2	10.09	10.06	10.20	9.4	23.7
4,4'-DDD	1	13.07	13.03	13.17	4.9	
	2	11.10	11.08	11.22	4.8	2.1

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY12

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477476

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608

ID: 0.53(mm)

GC Column(2): RTX-1701

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM TO	CONCENTRATION	%D
Endosulfan II	1	13.21	13.13 13.27	2.5	
	2	11.22	11.14 11.28	12	380.0
4,4'-DDT	1	13.79	13.78 13.92	13	
	2	11.49	11.47 11.61	3.2	306.3
Endrin aldehyde	1	14.04	13.96 14.10	2.6	
	2	12.12	12.04 12.18	21	707.7
Endosulfan sulfate	1	14.42	14.28 14.42	4.7	
	2	12.91	12.77 12.91	16	240.4
Methoxychlor	1	16.08	15.99 16.13	15	
	2	13.08	12.97 13.11	45	200.0
	1				
	2				
	1				
	2				
	1				
	2				

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY13

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477483 Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN04 Instrument ID (2): VARIAN05

GC Column(1): DB-608 ID: 0.53(mm) GC Column(2): RTX-1701 ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM TO	CONCENTRATION	%D
Heptachlor	1	7.86	7.80 7.90	1.1	
	2	6.11	6.08 6.18	0.14	685.7
Aldrin	1	8.66	8.62 8.72	0.33	
	2	6.74	6.64 6.74	0.41	24.2
Heptachlor epoxide	1	10.17	10.06 10.20	0.56	
	2	8.24	8.17 8.31	0.24	133.3
gamma-Chlordane	1	10.52	10.48 10.62	0.13	
	2	8.99	8.89 9.03	0.13	0.0
alpha-Chlordane	1	10.94	10.90 11.04	0.26	
	2	9.16	9.07 9.21	1.7	553.8
Dieldrin	1	11.82	11.74 11.88	4.5	
	2	9.63	9.59 9.73	0.46	878.3
Endrin	1	12.78	12.66 12.80	2.6	
	2	10.12	10.06 10.20	1.6	62.5
Endosulfan II	1	13.23	13.13 13.27	0.71	
	2	11.23	11.14 11.28	0.32	121.9

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY14DL

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477484 D20

Date(s) Analyzed: 02/13/92 02/13/92

Instrument ID (1): VARIAN05

Instrument ID (2): VARIAN04

GC Column(1): RTX-1701

ID: 0.53(mm)

GC Column(2): DB-608

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
gamma-BHC (Lindane)	1	5.66	5.61	5.71	34	
	2	7.04	6.98	7.08	46	35.3
Heptachlor	1	6.06	6.06	6.16	2.8	
	2	7.86	7.78	7.88	21	650.0
delta-BHC	1	7.83	7.73	7.83	5.5	
	2	8.12	8.12	8.22	14	154.5
Aldrin	1	6.72	6.63	6.73	12	
	2	8.63	8.60	8.70	3.2	275.0
Heptachlor epoxide	1	8.22	8.16	8.30	4.6	
	2	10.05	10.05	10.19	1.9	142.1
gamma-Chlordane	1	8.87	8.87	9.01	1.1	
	2	10.49	10.47	10.61	3.7	236.4
Endosulfan I	1	8.75	8.75	8.89	14	
	2	10.98	10.93	11.07	22	57.1
Dieldrin	1	9.59	9.57	9.71	3.3	
	2	11.84	11.71	11.85	5.9	78.8



10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY14DL

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477484 D20

Date(s) Analyzed: 02/13/92 02/13/92

Instrument ID (1): VARIAN05

Instrument ID (2): VARIAN04

GC Column(1): RTX-1701

ID: 0.53(mm)

GC Column(2): DB-608

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
Endrin	1	10.07	10.03	10.17	6.6	
	2	12.67	12.63	12.77	4.5	46.7
4,4'-DDD	1	11.14	11.07	11.21	3.8	
	2	13.06	13.02	13.16	5.4	42.1
4,4'-DDT	1	11.49	11.45	11.59	1.8	
	2	13.78	13.75	13.89	2.8	55.6
Endrin aldehyde	1	12.12	12.02	12.16	11	
	2	14.01	13.94	14.08	2.7	307.4
	1					
	2					
	1					
	2					
	1					
	2					
	1					
	2					

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY14

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477484

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608 ID: 0.53(mm) GC Column(2): RTX-1701 ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
alpha-BHC	1	6.00	6.00	6.10	0.36	
	2	4.88	4.79	4.89	2.8	677.8
gamma-BHC (Lindane)	1	7.04	7.00	7.10	49	
	2	5.65	5.63	5.73	40	22.5
Aldrin	1	8.64	8.62	8.72	4.3	
	2	6.72	6.64	6.74	15	248.8
Heptachlor epoxide	1	10.16	10.06	10.20	12	
	2	8.21	8.17	8.31	5.6	114.3
gamma-Chlordane	1	10.49	10.48	10.62	3.3	
	2	9.01	8.89	9.03	3.7	12.1
alpha-Chlordane	1	10.99	10.90	11.04	24	
	2	9.11	9.07	9.21	6.2	287.1
Endrin	1	12.69	12.66	12.80	4.4	
	2	10.07	10.06	10.20	14	218.2
4,4'-DDD	1	13.06	13.03	13.17	6.4	
	2	11.14	11.08	11.22	4.3	48.8

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY14

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477484

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608

ID: 0.53(mm)

GC Column(2): RTX-1701

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%D
=====	---	----	-----	-----	-----	-----
4,4'-DDT	1	13.79	13.78	13.92	1.5	
	2	11.49	11.47	11.61	2.4	60.0
Endrin aldehyde	1	14.00	13.96	14.10	3.1	
	2	12.11	12.04	12.18	12	287.1
Endosulfan sulfate	1	14.29	14.28	14.42	0.69	
	2	12.91	12.77	12.91	4.0	479.7
Methoxychlor	1	16.09	15.99	16.13	10.0	
	2	13.09	12.97	13.11	25	150.0
Endrin ketone	1	16.28	16.17	16.31	8.2	
	2	13.73	13.62	13.76	4.6	78.3
	1					
	2					
	1					
	2					
	1					
	2					

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY16

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477486

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN05

Instrument ID (2): VARIAN04

GC Column(1): RTX-1701

ID: 0.53(mm)

GC Column(2): DB-608

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
alpha-BHC	1	4.86	4.79	4.89	3.5	
	2	6.00	6.00	6.10	1.1	218.2
gamma-BHC (Lindane)	1	5.73	5.63	5.73	4.4	
	2	7.05	7.00	7.10	100	999.9
beta-BHC	1	7.35	7.27	7.37	6.6	
	2	7.21	7.21	7.31	3.9	69.2
Heptachlor	1	6.17	6.08	6.18	0.16	
	2	7.85	7.80	7.90	20	999.9
Aldrin	1	6.64	6.64	6.74	2.6	
	2	8.65	8.62	8.72	6.3	142.3
gamma-Chlordane	1	8.96	8.89	9.03	15	
	2	10.55	10.48	10.62	11	36.4
alpha-Chlordane	1	9.14	9.07	9.21	31	
	2	10.93	10.90	11.04	11	181.8
Endosulfan I	1	8.86	8.77	8.91	15	
	2	11.04	10.96	11.10	0.37	999.9

ORIGINAL  
(Red)

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY16

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477486

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN05

Instrument ID (2): VARIAN04

GC Column(1): RTX-1701 ID: 0.53(mm) GC Column(2): DB-608 ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
4,4'-DDE	1	9.35	9.31	9.45	42	100
	2	11.67	11.61	11.75	13	223.1
Dieldrin	1	9.63	9.59	9.73	37	
	2	11.80	11.74	11.88	74	100.0
Endrin	1	10.07	10.06	10.20	46	100
	2	12.70	12.66	12.80	160	247.8
Endosulfan II	1	11.22	11.14	11.28	190	
	2	13.21	13.13	13.27	64	196.9
Endrin aldehyde	1	12.11	12.04	12.18	360	100
	2	14.03	13.96	14.10	53	579.2
Endosulfan sulfate	1	12.91	12.77	12.91	190	100
	2	14.37	14.28	14.42	19	900.0
Methoxychlor	1	13.03	12.97	13.11	180	
	2	16.07	15.99	16.13	410	127.8
	1					
	2					

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY17

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477487

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN05

Instrument ID (2): VARIAN04

GC Column(1): RTX-1701 ID: 0.53(mm) GC Column(2): DB-608 ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
gamma-BHC (Lindane)	1	5.66	5.63	5.73	14	
	2	7.06	7.00	7.10	31	121.4
Heptachlor	1	6.12	6.08	6.18	21	
	2	7.84	7.80	7.90	16	31.2
Aldrin	1	6.69	6.64	6.74	15	
	2	8.66	8.62	8.72	14	7.1
Heptachlor epoxide	1	8.19	8.17	8.31	0.19	pu
	2	10.17	10.06	10.20	1.2	531.6
gamma-Chlordane	1	8.98	8.89	9.03	4.7	pu
	2	10.56	10.48	10.62	0.093	999.9
Dieldrin	1	9.65	9.59	9.73	31	
	2	11.81	11.74	11.88	47	51.6
Endrin	1	10.10	10.06	10.20	55	
	2	12.72	12.66	12.80	39	41.0
Endosulfan II	1	11.25	11.14	11.28	2.8	
	2	13.24	13.13	13.27	2.2	27.3

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY18

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477489 R1

Date(s) Analyzed: 02/18/92 02/18/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608

ID: 0.53(mm)

GC Column(2): RTX-1701

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM TO	CONCENTRATION	%D
Aldrin	1	8.61	8.60 8.70	0.36	
	2	6.71	6.63 6.73	0.76	111.1
Heptachlor epoxide	1	10.12	10.05 10.19	1.1	
	2	8.18	8.16 8.30	0.54	103.7
gamma-Chlordane	1	10.58	10.47 10.61	0.23	
	2	8.96	8.87 9.01	3.7	999.9
4,4'-DDE	1	11.64	11.59 11.73	2.8	
	2	9.34	9.30 9.44	3.2	14.3
Dieldrin	1	11.77	11.71 11.85	19	poor
	2	9.60	9.57 9.71	0.27	999.9
Endrin	1	12.73	12.63 12.77	2.6	per
	2	10.07	10.03 10.17	15	476.9
Endosulfan II	1	13.18	13.11 13.25	2.6	
	2	11.20	11.12 11.26	2.7	3.8
4,4'-DDT	1	13.80	13.75 13.89	0.72	
	2	11.47	11.45 11.59	0.48	50.0

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY18

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579H0 SDG No.: CHY11

Lab Sample ID : 477489

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608

ID: 0.53(mm)

GC Column(2): RTX-1701

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM TO	CONCENTRATION	%D
Heptachlor	1	7.85	7.80 7.90	0.30	
	2	6.08	6.08 6.18	0.12	150.0
Heptachlor epoxide	1	10.17	10.06 10.20	1.8	pos
	2	8.22	8.17 8.31	0.33	445.5
gamma-Chlordane	1	10.58	10.48 10.62	4.1	pos
	2	8.94	8.89 9.03	1.1	272.7
alpha-Chlordane	1	10.94	10.90 11.04	1.7	
	2	9.14	9.07 9.21	4.8	182.4
4,4'-DDE	1	11.67	11.61 11.75	1.3	
	2	9.36	9.31 9.45	2.5	92.3
Dieldrin	1	11.81	11.74 11.88	2.4	
	2	9.64	9.59 9.73	1.0	140.0
Endrin	1	12.71	12.66 12.80	4.2	
	2	10.10	10.06 10.20	5.0	19.0
4,4'-DDD	1	13.08	13.03 13.17	2.5	pos
	2	11.10	11.08 11.22	0.99	152.5



ORIGINAL  
(Red)

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY18

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477489

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608

ID: 0.53(mm)

GC Column(2): RTX-1701

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM TO	CONCENTRATION	%D
Endosulfan II	1	13.20	13.13 13.27	0.49	pool
	2	11.23	11.14 11.28	5.2	961.2
4,4'-DDT	1	13.79	13.78 13.92	6.1	
	2	11.50	11.47 11.61	2.2	177.3
Endrin aldehyde	1	14.02	13.96 14.10	1.0	pool
	2	12.12	12.04 12.18	7.5	650.0
Endosulfan sulfate	1	14.28	14.28 14.42	0.92	pool
	2	12.90	12.77 12.91	5.5	497.8
Methoxychlor	1	16.06	15.99 16.13	200	pool
	2	13.08	12.97 13.11	11	999.9
	1				
	2				
	1				
	2				
	1				
	2				

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY19

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477490

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608

ID: 0.53(mm)

GC Column(2): RTX-1701

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
alpha-BHC	1	6.02	6.00	6.10	0.081	
	2	4.85	4.79	4.89	0.15	85.2
Heptachlor	1	7.86	7.80	7.90	1.3	
	2	6.11	6.08	6.18	0.15	766.7
Aldrin	1	8.65	8.62	8.72	0.39	
	2	6.74	6.64	6.74	0.50	28.2
Heptachlor epoxide	1	10.17	10.06	10.20	0.41	
	2	8.21	8.17	8.31	0.63	53.7
4,4'-DDE	1	11.62	11.61	11.75	1.2	
	2	9.36	9.31	9.45	1.1	9.1
Dieldrin	1	11.82	11.74	11.88	6.3	
	2	9.64	9.59	9.73	0.10	999.9
Endrin	1	12.78	12.66	12.80	1.5	
	2	10.10	10.06	10.20	6.0	300.0
Endosulfan II	1	13.24	13.13	13.27	0.48	
	2	11.23	11.14	11.28	0.87	81.2

ORIGINAL  
Recd.

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY20

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477491

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN05

Instrument ID (2): VARIAN04

GC Column(1): RTX-1701

ID: 0.53(mm)

GC Column(2): DB-608

ID: 0.53(mm)

*Note: Chromatogram had continuous peaks similar Heights.*

ANALYTE	COL	RT	RT WINDOW FROM TO	CONCENTRATION	%D
beta-BHC	1	7.36	7.27 7.37	5.4	
	2	7.31	7.21 7.31	1.3	315.4
Heptachlor	1	6.08	6.08 6.18	1.8	
	2	7.86	7.80 7.90	17	844.4
Aldrin	1	6.72	6.64 6.74	8.1	
	2	8.65	8.62 8.72	5.6	44.6
Heptachlor epoxide	1	8.23	8.17 8.31	2.3	
	2	10.17	10.06 10.20	8.6	273.9
gamma-Chlordane	1	8.97	8.89 9.03	15	
	2	10.55	10.48 10.62	2.8	435.7
4,4'-DDE	1	9.35	9.31 9.45	11	
	2	11.68	11.61 11.75	11	0.0
Dieldrin	1	9.64	9.59 9.73	14	
	2	11.81	11.74 11.88	38	171.4
Endrin	1	10.08	10.06 10.20	34	
	2	12.71	12.66 12.80	52	52.9

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY20

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477491

Date(s) Analyzed: 02/07/92 02/07/92

Instrument ID (1): VARIAN05

Instrument ID (2): VARIAN04

GC Column(1): RTX-1701

ID: 0.53(mm)

GC Column(2): DB-608

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%D
=====	=====	=====	FROM	TO	=====	=====
Endosulfan II	1	11.23	11.14	11.28	73	
	2	13.21	13.13	13.27	20	265.0
Endrin aldehyde	1	12.11	12.04	12.18	190	
	2	14.03	13.96	14.10	17	999.9
Methoxychlor	1	13.03	12.97	13.11	120	
	2	16.08	15.99	16.13	250	108.3
	1					
	2					
	1					
	2					
	1					
	2					
	1					
	2					
	1					
	2					

10A

PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY21

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477492 R1

Date(s) Analyzed: 02/22/92 02/22/92

Instrument ID (1): VARIAN05

Instrument ID (2): VARIAN04

GC Column(1): RTX-1701

ID: 0.53(mm)

GC Column(2): DB-608

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
gamma-BHC (Lindane)	1	5.70	5.61	5.71	0.18	POOR
	2	7.07	6.98	7.08	54	999.9
Heptachlor	1	6.12	6.06	6.16	0.81	POOR
	2	7.87	7.78	7.88	6.8	739.5
Aldrin	1	6.66	6.63	6.73	1.4	
	2	8.66	8.60	8.70	2.0	42.9
Methoxychlor	1	13.04	12.95	13.09	110	
	2	16.05	15.97	16.11	180	63.6
gamma-Chlordane	1	8.99	8.87	9.01	14	
	2	10.55	10.47	10.61	5.7	145.6
alpha-Chlordane	1	9.16	9.06	9.20	16	POOR
	2	10.92	10.88	11.02	2.4	566.7
4,4'-DDE	1	9.37	9.30	9.44	6.4	
	2	11.67	11.59	11.73	6.5	1.6
Dieldrin	1	9.57	9.57	9.71	2.4	POOR
	2	11.80	11.71	11.85	30	999.9

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY22

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477493

Date(s) Analyzed: 02/08/92 02/08/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608 ID: 0.53(mm) GC Column(2): RTX-1701 ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW FROM	TO	CONCENTRATION	%D
gamma-Chlordane	1	10.54	10.48	10.62	2.1	999.9
	2	8.91	8.89	9.03	110	999.9
alpha-Chlordane	1	10.91	10.90	11.04	3.5	999.9
	2	9.21	9.07	9.21	17	385.7
4,4'-DDE	1	11.67	11.61	11.75	1.6	
	2	9.35	9.31	9.45	1.6	0.0
Dieldrin	1	11.81	11.74	11.88	9.4	
	2	9.65	9.59	9.73	6.2	51.6
4,4'-DDT	1	13.79	13.78	13.92	3.9	
	2	11.50	11.47	11.61	2.3	69.6
Methoxychlor	1	16.08	15.99	16.13	180	
	2	13.03	12.97	13.11	26	592.3
Endrin ketone	1	16.24	16.17	16.31	13	
	2	13.68	13.62	13.76	0.66	999.9
	1					
	2					

(Red)

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY23

Lab Name: COMPUCHEM.RTPContract: 68D10083Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11Lab Sample ID : 477494Date(s) Analyzed: 02/08/92 02/08/92Instrument ID (1): VARIAN04Instrument ID (2): VARIAN05GC Column(1): DB-608ID: 0.53(mm)GC Column(2): RTX-1701ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%D
			FROM	TO		
beta-BHC	1	7.21	7.21	7.31	0.87	
	2	7.35	7.27	7.37	1.8	106.9
Heptachlor	1	7.87	7.80	7.90	0.23	700
	2	6.18	6.08	6.18	1.2	421.7
Endrin	1	12.71	12.66	12.80	1.4	
	2	10.11	10.06	10.20	1.5	7.1
Endosulfan II	1	13.22	13.13	13.27	10.41	700
	2	11.23	11.14	11.28	1.6	290.2
Endrin aldehyde	1	13.99	13.96	14.10	0.68	700
	2	12.12	12.04	12.18	4.3	532.4
Methoxychlor	1	16.08	15.99	16.13	8.1	700
	2	13.09	12.97	13.11	1.4	478.6
	1					
	2					
	1					
	2					

10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY26

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477495

Date(s) Analyzed: 02/08/92 02/08/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608

ID: 0.53(mm)

GC Column(2): RTX-1701

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%D
=====	==	=====	=====	=====	=====	=====
gamma-BHC (Lindane)	1	7.05	7.00	7.10	3.6	
	2	5.69	5.63	5.73	0.091	999.9
beta-BHC	1	7.31	7.21	7.31	0.28	
	2	7.35	7.27	7.37	0.67	139.3
Heptachlor	1	7.86	7.80	7.90	1.9	
	2	6.10	6.08	6.18	0.19	900.0
Aldrin	1	8.65	8.62	8.72	0.58	
	2	6.72	6.64	6.74	0.56	3.6
4,4'-DDE	1	11.66	11.61	11.75	0.80	
	2	9.36	9.31	9.45	1.2	50.0
Dieldrin	1	11.82	11.74	11.88	3.3	
	2	9.63	9.59	9.73	0.26	999.9
Endrin	1	12.78	12.66	12.80	4.0	
	2	10.09	10.06	10.20	3.3	21.2
Endosulfan II	1	13.24	13.13	13.27	0.28	
	2	11.23	11.14	11.28	2.6	828.6



10A  
PESTICIDE IDENTIFICATION SUMMARY  
FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

CHY26

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Lab Sample ID : 477495

Date(s) Analyzed: 02/08/92 02/08/92

Instrument ID (1): VARIAN04

Instrument ID (2): VARIAN05

GC Column(1): DB-608

ID: 0.53(mm)

GC Column(2): RTX-1701

ID: 0.53(mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%D
=====	=====	=====	FROM	TO	=====	=====
4,4'-DDT	1	13.83	13.78	13.92	10	987.0
	2	11.50	11.47	11.61	0.92	
	1					
	2					
	1					
	2					
	1					
	2					
	1					
	2					
	1					
	2					
	1					
	2					
	1					
	2					

PROJECT NAME: Linfield Industrial Park  
Proj. NO.: 326309

EPA SITE NO.: PA091/2898  
REGION: 3

**SUPPORT DOCUMENTATION FOR THE REVIEW OF  
INORGANIC ANALYTICAL DATA PACKAGE**

CASE/SAS NO.: 17744  
TYPE OF ANALYSIS: low inorganic/CN-  
CONTRACT LABORATORY: Reg. TX  
APPLICABLE IFB OR SOW: 3/90  
REVIEWER: (b) (4)  
REVIEW DATE: May 2, 1992

APPLICABLE SAMPLE NO's.:  
MCJE02-26  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

THE FOLLOWING TABLE INDICATES  
AREAS WHICH WERE EXAMINED IN  
DETAIL, THE IDENTIFIED PROBLEM  
AREAS, AND SUPPORT DOCUMENTATION  
ATTACHMENTS:

	AREAS EXAMINED IN DETAIL					PROBLEM AREAS IDENTIFIED					SUPPORT DOCUMENTATION ATTACHMENTS				
	CHECK (✓) IF YES OR FOOTNOTE LETTER FOR COMMENTS BELOW					CHECK (✓) IF YES OR FOOTNOTE NUMBER FOR COMMENTS BELOW					CHECK (✓) IF YES OR IDENTIFY ATTACHMENT NO.				
	ALL APPLICABLE ANALYSES	ICP OR A.A. METALS	FURNACE METALS	COLD VAPOR MERCURY	CYANIDE	ALL APPLICABLE ANALYSES	ICP OR A.A. METALS	FURNACE METALS	COLD VAPOR MERCURY	CYANIDE	ALL APPLICABLE ANALYSES	ICP OR A.A. METALS	FURNACE METALS	COLD VAPOR MERCURY	CYANIDE
HOLDING TIMES	✓														
BLANK ANALYSIS RESULTS	✓														
MATRIX SPIKES (PRE-DIGESTION)	✓														
DUPLICATES	✓														
QUANTITATION OF RESULTS	✓														
DETECTION LIMITS/SENSITIVITY	✓														
INITIAL CALIBRATIONS	✓														
CONTINUING CALIBRATIONS	✓														
LABORATORY CONTROL STANDARDS	✓														
ICP LINEAR RANGE ANALYSIS	✓														
ICP INTERFERENCE CHECKS	✓														
ICP SERIAL DILUTIONS	✓														
GFAA POST-DIGESTION SPIKES	✓														
GFAA DUPLICATE BURNS	✓														
GFAA STANDARD ADDITIONS	✓														
OTHERS	✓														

COMMENTS: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

# SUMMARY OF BLANK ANALYSES

(Highest values for each category)

(Units = ug/l)

Case: \_\_\_\_\_

SDGs: \_\_\_\_\_

Metal	INITIAL CALIB. BLANKS			CONT. CALIB. BLANKS			PREP. BLANKS			FIELD BLANKS		
	AQ SDG#	FILT SDG#	SOL SDG#	AQ SDG#	FILT SDG#	SOL SDG#	AQ SDG#	FILT SDG#	SOL SDG#	AQ SPL#	FILT SPL#	SOL SPL#
	MC1802	MC	MC	MC1802	MC	MC	MC1802	MC	MC	MC1802	MC	
Al		X			X			X		(15.1)	N/A	N/A
As		X			X			X		(5.4)	N/A	N/A
Sb		X			X			X				
Ba		X			X			X				
Be		X			X			X				
Cd		X			X			X				
Ca				7.8						(29.1)		
Cr												
Co												
Cu												
Fe										(12.5)		
Pb												
Mg										(42.6)		
Mn												
Hg												
Ni												
K												
Se			2.2	2.3		2.5			(2.7)			
Ag												
Na												
Tl	2.1		2.4	4.0		(4.0)						
V												
Zn										(15.8)		

Highest values for each metal (unfiltered aqueous and solids) are circled and are applied to *all* samples. FILT field blank results apply *only* to filtered samples.

Comments: \_\_\_\_\_

**SUPPORT DOCUMENTATION**  
**for SDG MCJE02**

includes (~~aqueous~~ ~~solid~~) samples

**MCJE02-10, MCJE24**

U.S. EPA - CLP

2B

## CRDL STANDARD FOR AA AND ICP

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

AA CRDL Standard Source: EMSL-LV

ICP CRDL Standard Source: EMSL-LV

Concentration Units: ug/L

Analyte	CRDL Standard for AA			CRDL Standard for ICP				
	True	Found	%R	True	Initial Found	%R	Final Found	%R
Aluminum_								
Antimony_				120.0	124.40	103.7	123.33	102.8
Arsenic_	10.0	9.10	91.0					
Barium_								
Beryllium_				10.0	10.36	103.6	10.08	100.8
Cadmium_				10.0	10.28	102.8	10.28	102.8
Calcium_								
Chromium_				20.0	19.35	96.8	20.84	104.2
Cobalt_				100.0	103.49	103.5	100.28	100.3
Copper_				50.0	51.14	102.3	51.64	103.3
Iron_								
Lead_	3.0	2.40	80.0					
Magnesium_								
Manganese_				30.0	31.00	103.3	30.10	100.3
Mercury_								
Nickel_				80.0	85.81	107.3	70.36	87.9
Potassium_								
Selenium_	5.0	5.20	104.0					
Silver_				20.0	19.36	96.8	19.06	95.3
Sodium_								
Thallium_								
Vanadium_				100.0	101.70	101.7	101.70	101.7
Zinc_				40.0	45.63	114.1	44.94	112.3

U.S. EPA - CLP

3  
BLANKS

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Preparation Blank Matrix (soil/water): WATER

Preparation Blank Concentration Units (ug/L or mg/kg): ug/L

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum_	14.0	U	14.0	U	14.0	U	14.0	U	14.000	U	P
Antimony_	30.0	U	30.0	U	30.0	U	30.0	U	30.000	U	P
Arsenic__	2.0	U	2.0	U	2.0	U	2.0	U	2.000	U	F
Barium__	1.0	U	1.0	U	1.0	U	1.0	U	1.000	U	P
Beryllium	1.0	U	1.0	U	1.0	U	1.0	U	1.000	U	P
Cadmium__	3.0	U	3.0	U	3.0	U	3.0	U	3.000	U	P
Calcium__	7.0	U	7.0	U	7.0	U	7.0	U	7.000	U	P
Cromium_	3.0	U	3.0	U	3.0	U	3.0	U	3.000	U	P
Cobalt__	4.0	U	4.0	U	4.0	U	4.0	U	4.000	U	P
Copper__	3.0	U	3.0	U	3.0	U	3.0	U	3.000	U	P
Iron_____	6.0	U	6.0	U	6.0	U	6.0	U	6.000	U	P
Lead_____	1.0	U	1.0	U	1.0	U	-1.0	B	1.000	U	F
Magnesium	37.0	U	37.0	U	37.0	U	37.0	U	37.000	U	P
Manganese	2.0	U	2.0	U	2.0	U	2.0	U	2.000	U	P
Mercury__	0.2	U	0.2	U	0.2	U	0.2	U	0.200	U	CV
Nickel__	22.0	U	22.0	U	22.0	U	22.0	U	22.000	U	P
Potassium	722.0	U	722.0	U	722.0	U	722.0	U	722.000	U	P
Selenium_	2.0	U	2.0	U	2.0	U	2.3	B	2.000	U	F
Silver__	2.0	U	2.0	U	2.0	U	2.0	U	2.000	U	P
Sodium__	30.0	U	30.0	U	30.0	U	30.0	U	30.000	U	P
Thallium_	2.1	B	2.0	U	2.0	U	4.0	B	2.000	U	F
Vanadium_	4.0	U	4.0	U	4.0	U	4.0	U	4.000	U	P
Zinc_____	2.0	U	2.0	U	2.0	U	2.0	U	2.000	U	P
Cyanide__	10.0	U	10.0	U	10.0	U	10.0	U	10.000	U	AS

FORM III - IN

3/90

000021

## U.S. EPA - CLP

3  
BLANKS

b Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Preparation Blank Matrix (soil/water):

Preparation Blank Concentration Units (ug/L or mg/kg):

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum_			14.0	U							P
Antimony_			30.0	U							P
Arsenic__			2.0	U	2.3	B	2.0	U			F
Barium__			1.0	U							P
Beryllium			1.0	U							P
Cadmium__			3.0	U							P
Calcium__			7.8	B							P
Cesium__			3.0	U							P
Cobalt__			4.0	U							P
Copper__			3.0	U							P
Iron_____			6.0	U							P
Lead_____			-1.2	B	-1.5	B	-1.5	B			F
Magnesium			37.0	U							P
Manganese			2.0	U							P
Mercury__											
Nickel__			22.0	U							P
Potassium			722.0	U							P
Selenium_			2.0	U	2.0	U	2.0	U			F
Silver__			2.0	U							P
Sodium__			30.0	U							P
Thallium_			2.0	U	2.0	U	2.0	U			F
Vanadium_			4.0	U							P
Zinc_____			2.0	U							P
Cyanide__											

U.S. EPA - CLP

3  
BLANKS

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Preparation Blank Matrix (soil/water):

Preparation Blank Concentration Units (ug/L or mg/kg):

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum_											
Antimony_											
Arsenic_			2.0	U							F
Barium_											
Beryllium_											
Cadmium_											
Calcium_											
Chromium_											
Cobalt_											
Copper_											
Iron_											
Lead_			-1.7	B							F
Magnesium_											
Manganese_											
Mercury_											
Nickel_											
Potassium_											
Selenium_			2.8	B	2.0	U	2.0	U			F
Silver_											
Sodium_											
Thallium_											
Vanadium_											
Zinc_											
Cyanide_											



4  
ICP INTERFERENCE CHECK SAMPLE

Lab Name: KEYSTONE LAB-HOUSTON      Contract: 68-D0-0147  
Lab Code: KEYTX      Case No.: 17744      SAS No.:      SDG No.: MCJE02  
ICP ID Number: TJA61      ICS Source: EPALV90

Concentration Units: ug/L

Analyte	True		Initial Found			Final Found		
	Sol. A	Sol. AB	Sol. A	Sol. AB	%R	Sol. A	Sol. AB	%R
Aluminum_	540055	536472	496982	498526.8	92.9	490764	500246.4	93.2
Antimony_			22	-49.4		25	7.5	
Arsenic_								
Barium_		502	12	456.3	90.9	11	456.1	90.9
Beryllium		480	1	472.6	98.5	1	474.2	98.8
Cadmium_		907	-2	868.8	95.8	-4	873.0	96.2
Calcium_	494040	512228	503913	505909.1	98.8	501572	508606.9	99.3
Chromium_	21	529	16	493.2	93.2	17	493.2	93.2
Cobalt_		477	11	441.0	92.4	13	439.7	92.2
Copper_		543	-20	528.8	97.4	-21	531.3	97.8
Cu_	206236	199845	188128	188247.7	94.2	186432	188684.2	94.4
Lead_								
Magnesium	531358	527530	515568	517732.0	98.1	509709	519827.5	98.5
Manganese	34	496	-33	451.2	91.0	-34	454.3	91.6
Mercury_								
Nickel_		940	23	858.5	91.3	19	851.2	90.6
Potassium			-1312	-989.2		-1323	-1290.3	
Selenium_								
Silver_		960	-38	859.8	89.6	-36	863.3	89.9
Sodium_			1764	1767.7		1730	1760.6	
Thallium_								
Vanadium_		509	33	481.3	94.6	33	486.9	95.6
Zinc_	216	1208	218	1160.2	96.0	218	1165.7	96.5

U.S. EPA - CLP

SA  
SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

MCJE07S

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix: WATER

Level (low/med): LOW

% Solids for Sample: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

Analyte	Control Limit %R	Spiked Sample Result (SSR)	C	Sample Result (SR)	C	Spike Added (SA)	%R	Q	M
Aluminum	75-125	2681.0000		1089.6200		2000.00	79.6	P	
Antimony	75-125	447.9800		30.0000	U	500.00	89.6	P	
Arsenic	75-125	50.4000		7.2000	B	40.00	108.0	F	
Barium	75-125	1764.9399		65.3500	B	2000.00	85.0	P	
Beryllium	75-125	47.1700		1.0000	U	50.00	94.3	P	
Cadmium	75-125	46.6200		3.0000	U	50.00	93.2	P	
Calcium								NR	
Chromium	75-125	177.7200		3.0000	U	200.00	88.9	P	
Cobalt	75-125	442.8800		4.0000	U	500.00	88.6	P	
Copper	75-125	226.9300		4.9600	B	250.00	88.8	P	
Iron	75-125	2252.9500		930.4400	SR & SA	1000.00	132.20	OKN	P
Lead	75-125	19.7000		1.1000	B	20.00	93.0	F	
Magnesium								NR	
Manganese	75-125	495.6600		51.6700		500.00	88.8	P	
Mercury	75-125	1.0000		0.2000	U	1.00	100.0	CV	
Nickel	75-125	441.9000		22.0000	U	500.00	88.4	P	
Potassium								NR	
Selenium	75-125	10.6000		2.0000	U	10.00	106.0	F	
Silver	75-125	41.9400		2.0000	U	50.00	83.9	P	
Sodium								NR	
Thallium	75-125	50.8000		2.0000	U	50.00	101.6	F	
Vanadium	75-125	442.8300		4.0000	U	500.00	88.6	P	
Zinc	75-125	477.7300		27.6800		500.00	90.0	P	
Cyanide	75-125	97.1700		10.0000	U	100.00	97.2	AS	

Comments:

Metals Se, AsIDLs 2.2 ug/L

U = &lt;IDL

Graphite Furnace Spike Recovery Evaluation Form

17744

SP6MCTE02

Sample ID	Instr. Level Result ug/L	PDS Recovery %	Diluted Result ug/L	Diluted PDS Recovery %	MSA Result (if needed) ug/L	Final Result Reported ug/L
MCTE02	6.2	55			2.0 <sup>low</sup>	2.0 ✓
03	ND	105				2.0 ✓
04	2.8	55			ND	2.0 ✓
05	2.0	94				2.0 ✓
06	2.0	107				2.0 ✓
07	ND	108				2.0 ✓
07D	ND	94				2.0 ✓
07S	10.6	—				10.6 %
08	ND	93				2.0
09	2.8	81	<del>ND</del> 2.5	86		2.5 ✓
10	3.3	78				
24	ND	87				
MCTE02	10.8	106				10.8 ✓
03	5.6	132	<del>ND</del> 8.7	86		8.1 ✓
04	5.8	107				5.8 ✓
05	6.8	101				6.8 ✓
06	7.2	105				7.2 ✓
07	7.2	94				7.2 ✓
07D	6.6	94				6.6 ✓
07S	50.4	—				108 %
08	8.9	104				8.9 ✓
09	6.7	107				6.7 ✓
10	6.4	89				6.4
24	5.4	82				5.4 ✓

Se

As

Metals Pb, Tl  
 IDLs 1.2 ug/l  
 U = < IDL

17744  
 SP6MCJF02

Graphite Furnace Spike Recovery Evaluation Form

Sample ID	Instr. Level Result ug/l	PDS Recovery %	Diluted Result ug/l	Diluted PDS Recovery %	MSA Result (if needed) ug/l	Final Result Reported ug/l	100
MCJF02	ND	94				1 U	
03	ND	104				↓	
04	ND	108				↓	
05	ND	120				↓	
06	22	119.			9.5x2	19.0	(23.5)
07	1.1	106				1.1	✓
07D	1.5	111				1.5	✓
07S	19.7	-				9.3 %	(85%)
08	59.9	97				59.9	71.5
09	N.D	100				1 U	
10	2.7	105				2.7	✓
24	ND	98				1 U	
MCJF02	ND	105				2 U	
03	ND	104				↓	
04	ND	108				↓	
05	ND	104				↓	
06	ND	105				↓	
07	ND	100				↓	
07D	ND	109				↓	
07S	50.8	-				102 %	
08	ND	108				2 U	
09	ND	105				↓	
10	ND	111				↓	
24	ND	105				↓	

Pb ICP IDL = 25 ug/l

ORIGINAL  
(Red)

U.S. EPA - CLP

KEYSTONE D.C.# 17744-3- 9

5B  
POST DIGEST SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

MCJE07A

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix (soil/water): WATER

Level (low/med): LOW

Concentration Units: ug/L

Analyte	Control Limit %R	Spiked Sample Result (SSR) C	Sample Result (SR) C	Spike Added (SA)	%R	Q/M
Aluminum_						NR
Antimony_						NR
Arsenic__						NR
Barium___						NR
Beryllium						NR
Cadmium__						NR
Calcium__						NR
Chromium_						NR
Cobalt___						NR
Copper___						NR
Iron_____		2885.46	930.44	1900.0	102.9	P
Lead_____						NR
Magnesium						NR
Manganese						NR
Mercury__						NR
Nickel___						NR
Potassium						NR
Selenium_						NR
Silver___						NR
Sodium___						NR
Thallium_						NR
Vanadium_						NR
Zinc_____						NR
Cyanide__						NR

Comments:

FORM V (PART 2) - IN

Rev. 3/90

000026

## U.S. EPA - CLP

6  
DUPLICATES

EPA SAMPLE NO.

MCJE07D

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix (soil/water): WATER

Level (low/med): LOW

% Solids for Sample: 0.0

% Solids for Duplicate: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

Analyte	Control Limit	Sample (S)	C	Duplicate (D)	C	RPD	Q	M
Aluminum		1089.6200		1115.5200		2.3		P
Antimony		30.0000	U	30.0000	U			P
Arsenic		7.2000	B	6.6000	B	8.7		F
Barium		65.3500	B	66.0400	B	1.1		P
Beryllium		1.0000	U	1.0000	U			P
Cadmium		3.0000	U	3.0000	U			P
Calcium	5000.0	14513.1201		14709.2402		1.3		P
Chromium		3.0000	U	3.0000	U			P
Cobalt		4.0000	U	4.0000	U			P
Copper		4.9600	B	4.4600	B	10.6		P
Iron		930.4400		926.0000		0.5		P
Lead		1.1000	B	1.5000	B	30.8		F
Magnesium	5000.0	5787.0801		5925.9600		2.4		P
Manganese	15.0	51.6700		52.1200		0.9		P
Mercury		0.2000	U	0.2000	U			CV
Nickel		22.0000	U	22.0000	U			P
Potassium		2268.8101	B	2322.5701	B	2.3		P
Selenium		2.0000	U	2.0000	U			F
Silver		2.0000	U	2.0000	U			P
Sodium	5000.0	6152.1001		6300.7798		2.4		P
Thallium		2.0000	U	2.0000	U			F
Vanadium		4.0000	U	4.0000	U			P
Zinc	20.0	27.6800		27.6300		0.2		P
Cyanide		10.0000	U	10.0000	U			AS

ORIGINAL  
(Red)

KEYSTONE D.C.# 17744-3- 12

U.S. EPA - CLP

8  
STANDARD ADDITION RESULTS

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Concentration Units: ug/L

Lab reported as  
2.0 not 2.0 ug/L

EPA Sample No.	An	0 ADD ABS	1 ADD CON ABS	2 ADD CON ABS	3 ADD CON ABS	Final Conc.	r	Q
MCJE06	PB	0.031	5.00 0.047	10.00 0.062	15.00 0.080	9.5	0.9993	
MCJE02	SE	0.017	5.00 0.027	10.00 0.052	15.00 0.091	2.0	0.9672	+
MCJE02	SE	0.007	5.00 0.025	10.00 0.050	15.00 0.041	4.6	0.8670	+
MCJE03	SE	0.005	5.00 0.040	10.00 0.070	15.00 0.115	0.5	0.9965	+
MCJE03	SE	0.005	5.00 0.020	10.00 0.043	15.00 0.065	0.7	0.9961	+
MCJE04	SE	0.003	5.00 0.038	10.00 0.055	15.00 0.138	0.5	0.9535	+
MCJE04	SE	0.011	5.00 0.024	10.00 0.054	15.00 0.054	3.7	0.9446	+

0.007 suggests  
sample is ND  
not 4.6

R same

MCJE02, 03, and 04  
are potable well samples  
and should not have  
serious matrix effects.  
This, coupled with  
erratic response, could  
suggest analytical  
malfunction or  
mis-spiking.

FORM VIII - IN

3/90

000029

U.S. EPA - CLP

9  
ICP SERIAL DILUTIONS

EPA SAMPLE NO.  
MCJE07L

Lab Name: KEYSTONE LAB-HOUSTON Contract: 68-DO-0147  
Lab Code: KEYTX Case No.: 17744 SAS No.: SDG No.: MCJE02  
Matrix (soil/water): WATER Level (low/med): LOW

Concentration Units: ug/L

Analyte	Initial Sample Result (I)	C	Serial Dilution Result (S)	C	% Difference	Q/M
Aluminum	1089.62		822.15	B	24.5	E P
Antimony	30.00	U	150.00	U		P
Arsenic						NR
Barium	65.35	B	65.70	B	0.5	P
Beryllium	1.00	U	5.00	U		P
Cadmium	3.00	U	15.00	U		P
Calcium	14513.12		15280.20	B	5.3	P
Chromium	3.00	U	15.00	U		P
Cobalt	4.00	U	20.00	U		P
Copper	4.96	B	15.00	U	100.0	P
Iron	930.44		767.70		17.5	E P
Lead						NR
Magnesium	5787.08		6042.40	B	4.4	P
Manganese	51.67		51.65	B	0.0	P
Mercury						NR
Nickel	22.00	U	110.00	U		P
* Potassium	2268.81	B	4408.55	B	94.3	P ok *
Selenium						NR
Silver	2.00	U	10.00	U		P
Sodium	6152.10		6433.75	B	4.6	P
Thallium						NR
Vanadium	4.00	U	20.00	U		P
Zinc	27.68		75.95	B	174.4	P ok both are (B)

$$K' \left( \frac{S-I}{S} \right) < 10L, \text{ no impact}$$



ORIGINAL  
(Red)

KEYSTONE D.C.# 17744-3-14

U.S. EPA - CLP

10

Instrument Detection Limits (Quarterly)

Lab Name: KEYSTONE LAB-HOUSTON Contract: 68-D0-0147  
Lab Code: KEYTX Case No.: 17744 SAS No.: SDG No.: MCJE02  
ICP ID Number: Date: 02/18/92  
Flame AA ID Number:  
Furnace AA ID Number: PE3030Z

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum			200		NR
Antimony			60		NR
Arsenic			10		NR
Barium			200		NR
Beryllium			5		NR
Cadmium			5		NR
Calcium			5000		NR
Chromium			10		NR
Cobalt			50		NR
Copper			25		NR
Iron			100		NR
Lead			3		NR
Magnesium			5000		NR
Manganese			15		NR
Mercury			0.2		NR
Nickel			40		NR
Potassium			5000		NR
Selenium			5		NR
Silver			10		NR
Sodium			5000		NR
Thallium	276.80	BZ	10	2.0	F
Vanadium			50		NR
Zinc			20		NR

Comments:

FORM X - IN

000031

U.S. EPA - CLP

10

Instrument Detection Limits (Quarterly)

ORIGINAL  
JEC

Lab Name: KEYSTONE LAB-HOUSTON Contract: 68-D0-0147  
Lab Code: KEYTX Case No.: 17744 SAS No.: SDG No.: MCJE02  
ICP ID Number: TJA61 Date: 02/10/92  
Flame AA ID Number:  
Furnace AA ID Number:

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum	308.20		200	14.0	P
Antimony	206.80		60	30.0	P
Arsenic			10		NR
Barium	493.40		200	1.0	P
Beryllium	313.00		5	1.0	P
Cadmium	228.80		5	3.0	P
Calcium	317.80		5000	7.0	P
Chromium	267.70		10	3.0	P
Cobalt	228.60		50	4.0	P
Copper	324.70		25	3.0	P
Iron	259.90		100	6.0	P
Lead			3		NR
Magnesium	279.00		5000	37.0	P
Manganese	257.60		15	2.0	P
Mercury			0.2		NR
Nickel	231.60		40	22.0	P
Potassium	766.40		5000	722.0	P
Selenium			5		NR
Silver	328.00		10	2.0	P
Sodium	588.90		5000	30.0	P
Thallium			10		NR
Vanadium	292.40		50	4.0	P
Zinc	213.80		20	2.0	P

Comments:

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FORM X - IN

000032

ORIGINAL  
(Red)

U.S. EPA - CLP

10

Instrument Detection Limits (Quarterly)

Lab Name: KEYSTONE LAB-HOUSTON Contract: 68-D0-0147  
Lab Code: KEYTX Case No.: 17744 SAS No.: SDG No.: MCJE02  
ICP ID Number: Date: 02/14/92  
Flame AA ID Number:  
Furnace AA ID Number: PE560

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum	283.30	BD	200	1.0	NR
Antimony			60		NR
Arsenic			10		NR
Barium			200		NR
Beryllium			5		NR
Cadmium			5		NR
Calcium			5000		NR
Chromium			10		NR
Cobalt			50		NR
Copper			25		NR
Iron			100		NR
Lead			3		F
Magnesium			5000		NR
Manganese			15		NR
Mercury			0.2		NR
Nickel			40		NR
Potassium			5000		NR
Selenium			5		NR
Silver			10		NR
Sodium			5000		NR
Thallium			10		NR
Vanadium			50		NR
Zinc			20		NR

Comments:

FORM X - IN

000033

## U.S. EPA - CLP

10

## Instrument Detection Limits (Quarterly)

10/15/92  
10/15/92

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

ICP ID Number:

Date: 01/15/92

Flame AA ID Number: MASS0A

Furnace AA ID Number:

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum_			200		NR
Antimony_			60		NR
Arsenic__			10		NR
Barium___			200		NR
Beryllium			5		NR
Cadmium__			5		NR
Calcium__			5000		NR
Chromium_			10		NR
Cobalt___			50		NR
Copper___			25		NR
Iron_____			100		NR
Lead_____			3		NR
Magnesium			5000		NR
Manganese			15		NR
Mercury__	253.70		0.2	0.2	CV
Nickel___			40		NR
Potassium			5000		NR
Selenium_			5		NR
Silver___			10		NR
Sodium___			5000		NR
Thallium_			10		NR
Vanadium_			50		NR
Zinc_____			20		NR

Comments:

FORM X - IN

000034

U.S. EPA - CLP

10

## Instrument Detection Limits (Quarterly)

Lab Name: KEYSTONE LAB-HOUSTON Contract: 68-DO-0147  
 Lab Code: KEYTX Case No.: 17744 SAS No.: SDG No.: MCJE02  
 ICP ID Number: Date: 02/22/92  
 Flame AA ID Number:  
 Furnace AA ID Number: TJASH21

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum	197.30	BS	200	2.0	NR
Antimony			60		NR
Arsenic			10		F
Barium			200		NR
Beryllium			5		NR
Cadmium			5		NR
Calcium			5000		NR
Chromium			10		NR
Cobalt			50		NR
Copper			25		NR
Iron	196.00	BS	100	2.0	NR
Lead			3		NR
Magnesium			5000		NR
Manganese			15		NR
Mercury			0.2		NR
Nickel			40		NR
Potassium			5000		NR
Selenium			5		F
Silver			10		NR
Sodium			5000		NR
Thallium			10		NR
Vanadium			50		NR
Zinc			20		NR

Comments:

**SUPPORT DOCUMENTATION**

**for SDG MCJE 11**

includes (~~aqueous~~, solid) samples

**MCJE-11-23, 26**

## U.S. EPA - CLP

2B

## CRDL STANDARD FOR AA AND ICP

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

AA CRDL Standard Source: EMSL-LV

ICP CRDL Standard Source: EMSL-LV

Concentration Units: ug/L

Analyte	CRDL Standard for AA			True	CRDL Standard for ICP			
	True	Found	%R		Initial Found	%R	Final Found	%R
Aluminum_								
Antimony_				120.0	122.02	101.7	121.07	100.9
Arsenic_	10.0	9.30	93.0					
Barium_								
Beryllium				10.0	9.98	99.8	9.74	97.4
Cadmium_				10.0	8.88	88.8	10.78	107.8
Calcium_								
Chromium_				20.0	20.26	101.3	20.26	101.3
Cobalt_				100.0	98.26	98.3	95.35	95.3
Copper_				50.0	51.14	102.3	48.77	97.5
Iron_								
Lead_	3.0	3.10	103.3					
Magnesium								
Manganese				30.0	29.28	97.6	28.86	96.2
Mercury_								
Nickel_				80.0	76.18	95.2	81.63	102.0
Potassium								
Selenium_	5.0	6.40	128.0					
Silver_				20.0	18.58	92.9	16.89	84.4
Sodium_								
Thallium_	10.0	11.90	119.0					
Vanadium_				100.0	98.64	98.6	95.58	95.6
Zinc_				40.0	43.11	107.8	42.57	106.4

## U.S. EPA - CLP

3  
BLANKS

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Preparation Blank Matrix (soil/water): SOIL

Preparation Blank Concentration Units (ug/L or mg/kg): MG/KG

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum_	14.0	U	14.0	U	14.0	U	14.0	U	2.800	U	P
Antimony_	30.0	U	30.0	U	30.0	U	30.0	U	6.000	U	P
Arsenic_	2.0	U	2.0	U	2.0	U	2.0	U	0.400	U	F
Barium_	1.0	U	1.0	U	1.0	U	1.0	U	0.200	U	P
Beryllium	1.0	U	1.0	U	1.0	U	1.0	U	0.200	U	P
Cadmium_	3.0	U	3.0	U	3.0	U	3.0	U	0.600	U	P
Calcium_	7.0	U	7.0	U	7.0	U	7.0	U	1.400	U	P
Cesium_	3.0	U	3.0	U	3.0	U	3.0	U	0.600	U	P
Cobalt_	4.0	U	4.0	U	4.0	U	4.0	U	0.800	U	P
Copper_	3.0	U	3.0	U	3.0	U	3.0	U	0.600	U	P
Iron_	6.0	U	6.0	U	6.0	U	6.0	U	1.200	U	P
Lead_	1.0	U	1.0	U	1.0	U	-1.0	B	0.200	U	F
Magnesium	37.0	U	37.0	U	37.0	U	37.0	U	7.400	U	P
Manganese	2.0	U	2.0	U	2.0	U	2.0	U	0.400	U	P
Mercury_	0.2	U	0.2	U	0.2	U	0.2	U	0.100	U	CV
Nickel_	22.0	U	22.0	U	22.0	U	22.0	U	4.400	U	P
Potassium	722.0	U	722.0	U	722.0	U	722.0	U	144.400	U	P
Selenium_	2.2	B✓	2.4	B✓	2.0	U	2.5	B✓	0.540	B	F✓
Silver_	2.0	U	2.0	U	2.0	U	2.0	U	0.400	U	P
Sodium_	30.0	U✓	30.0	U	30.0	U	30.0	U	6.000	U	P
Thallium_	2.1	B✓	2.0	U	2.0	U	4.0	B✓	0.400	U	F
Vanadium_	4.0	U	4.0	U	4.0	U	4.0	U	0.800	U	P
Zinc_	2.0	U	2.0	U	2.0	U	2.0	U	0.400	U	P
Cyanide_	10.0	U	10.0	U	10.0	U	10.0	U	0.500	U	AS



ORIGINAL  
(Red)

U.S. EPA - CLP

3  
BLANKS

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Preparation Blank Matrix (soil/water):

Preparation Blank Concentration Units (ug/L or mg/kg):

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum_			14.0	U							P
Antimony_			30.0	U							P
Arsenic_			2.0	U	2.3	B	2.0	U			F
Barium_			1.0	U							P
Beryllium_			1.0	U							P
Cadmium_			3.0	U							P
Calcium_			7.0	U							P
Chromium_			3.0	U							P
Cobalt_			4.0	U							P
Copper_			3.0	U							P
Iron_			6.0	U							P
Lead_			1.0	U	1.0	U	1.0	U			F
Magnesium_			37.0	U							P
Manganese_			2.0	U							P
Mercury_											
Nickel_			22.0	U							P
Potassium_			722.0	U							P
Selenium_			2.0	U	2.0	B	2.0	U			F
Silver_			2.0	U							P
Sodium_			30.0	U							P
Thallium_			2.0	U	2.0	U	2.0	U			F
Vanadium_			4.0	U							P
Zinc_			2.0	U							P
Cyanide_											

## U.S. EPA - CLP

3  
BLANKS

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DQ-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Preparation Blank Matrix (soil/water):

Preparation Blank Concentration Units (ug/L or mg/kg):

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum											
Antimony											
Arsenic			2.0	U	2.0	U	-2.2	B			F
Barium											
Beryllium											
Cadmium											
Calcium											
Cesium											
Cobalt											
Copper											
Iron											
Lead			1.0	U	-1.5	B	-1.3	B			F
Magnesium											
Manganese											
Mercury											
Nickel											
Potassium											
Selenium			2.0	U	2.8	B	2.0	U			F
Silver											
Sodium											
Thallium			2.0	U	2.0	U	2.0	U			F
Vanadium											
Zinc											
Cyanide											

ORIGINAL

(Red)

## U.S. EPA - CLP

3  
BLANKS

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Preparation Blank Matrix (soil/water):

Preparation Blank Concentration Units (ug/L or mg/kg):

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum_											
Antimony_											
Arsenic_			2.0	U	2.0	U	2.0	U			F
Barium_											
Beryllium_											
Cadmium_											
Calcium_											
Chromium_											
Cobalt_											
Copper_											
Iron_											
Lead_			1.0	U	-1.1	B					F
Magnesium_											
Manganese_											
Mercury_											
Nickel_											
Potassium_											
Selenium_			2.0	U	2.0	U	2.0	U			F
Silver_											
Sodium_											
Thallium_			2.0	U	2.0	U	2.0	U			F
Vanadium_											
Zinc_											
Cyanide_											

FORM III - IN

3/90

000034

## U.S. EPA - CLP

3  
BLANKS

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Preparation Blank Matrix (soil/water):

Preparation Blank Concentration Units (ug/L or mg/kg):

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						C	Prepa- ration Blank	C	M
			1	C	2	C	3	C				
Aluminum_												
Antimony_												
Arsenic_												
Barium_												
Beryllium_												
Cadmium_												
Calcium_												
Cromium_												
Cobalt_												
Copper_												
Iron_												
Lead_												
Magnesium_												
Manganese_												
Mercury_												
Nickel_												
Potassium_												
Selenium_												
Silver_												
Sodium_												
Thallium_			2.0	U	2.0	U	2.0	U				F
Vanadium_												
Zinc_												
Cyanide_												

FORM III - IN

3/90

000035

ORIGINAL  
(Red)

U.S. EPA - CLP

3  
BLANKS

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Preparation Blank Matrix (soil/water):

Preparation Blank Concentration Units (ug/L or mg/kg):

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum_											
Antimony_											
Arsenic_											
Barium_											
Beryllium_											
Cadmium_											
Calcium_											
Chromium_											
Cobalt_											
Copper_											
Iron_											
Lead_											
Magnesium_											
Manganese_											
Mercury_											
Nickel_											
Potassium_											
Selenium_											
Silver_											
Sodium_											
Thallium_			2.0	U	2.0	U	2.0	U			F
Vanadium_											
Zinc_											
Cyanide_											

FORM III - IN

3/90

000036

ORIGINAL  
(Red)

KEYSTONE D.C.# 17744-3-6

U.S. EPA - CLP

3  
BLANKS

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Preparation Blank Matrix (soil/water):

Preparation Blank Concentration Units (ug/L or mg/kg):

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L)						Prepa- ration Blank	C	M
			1	C	2	C	3	C			
Aluminum											
Antimony											
Arsenic											
Barium											
Beryllium											
Cadmium											
Calcium											
Chromium											
Cobalt											
Copper											
Iron											
Lead											
Magnesium											
Manganese											
Mercury											
Nickel											
Potassium											
Selenium											
Silver											
Sodium											
Thallium			2.0	U	2.0	U	2.0	U			F
Vanadium											
Zinc											
Cyanide											

ORIGINAL  
(Red)

KEYSTONE D.C.# 17744-3- 7

U.S. EPA - CLP

4

ICP INTERFERENCE CHECK SAMPLE

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

ICP ID Number: TJA61

ICS Source: EPALV90

Concentration Units: ug/L

Analyte	True		Initial Found			Final Found		
	Sol. A	Sol. AB	Sol. A	Sol. AB	%R	Sol. A	Sol. AB	%R
Aluminum_	540055	536472	490703	498407.6	92.9	492685	489829.9	91.3
Antimony_			14	7.0		29	-14.8	
Arsenic_								
Barium_		502	12	453.8	90.4	12	439.1	87.5
Beryllium_		480	0	449.4	93.6	0	432.6	90.1
Cadmium_		907	-2	867.8	95.7	-2	857.1	94.5
Calcium_	494040	512228	490112	497994.1	97.2	488999	485989.6	94.9
Chromium_	21	529	19	482.0	91.1	17	468.5	88.6
Cobalt_		477	11	431.7	90.5	12	419.8	88.0
Copper_		543	-21	516.7	95.1	-21	502.0	92.4
Copper_	206236	199845	181855	185392.4	92.8	179605	179494.0	89.8
Lead_								
Magnesium_	531358	527530	501252	510837.4	96.8	508379	506125.0	95.9
Manganese_	34	496	-30	446.2	90.0	-28	433.4	87.4
Mercury_								
Nickel_		940	18	833.2	88.6	2	792.0	84.3
Potassium_			-1242	-1241.7		-675	-1082.7	
Selenium_								
Silver_		960	-38	869.1	90.5	-38	863.7	90.0
Sodium_			1636	1694.6		1687	1684.4	
Thallium_								
Vanadium_		509	35	476.8	93.7	38	468.4	92.0
Zinc_	216	1208	212	1146.4	94.9	216	1135.3	94.0

FORM IV - IN

3/90

000038

## U.S. EPA - CLP

5A  
SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

MCJE18S

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix: SOIL

Level (low/med): LOW

% Solids for Sample: 76.0

Concentration Units (ug/L or mg/kg dry weight): MG/KG

Analyte	Control Limit %R	Spiked Sample Result (SSR)	C	Sample Result (SR)	C	Spike Added (SA)	%R	Q	M
Aluminum									NR
Antimony	75-125	121.7868		7.8947	U	131.58	92.6		IF
Arsenic	75-125	16.8684		6.6579		10.53	97.0		IF
Barium	75-125	531.6974		52.7026		526.32	91.0		IF
Beryllium	75-125	13.0105		0.5105	B	13.16	95.0		IF
Cadmium	75-125	13.0158		0.7895	U	13.16	98.9		IF
Calcium									NR
Chromium	75-125	63.6000		10.6316		52.63	100.6		IF
Chloride	75-125	128.4421		6.3658	B	131.58	92.8		IF
Copper	75-125	73.5263		9.9684		65.79	96.6		IF
Iron									NR
Lead		239.4737		181.5789		5.26	1100.7		IF
Magnesium									NR
Manganese	75-125	429.7684		317.6579		131.58	85.2		IF
Mercury	75-125	0.6579		0.1316	U	0.66	99.7		CV
Nickel	75-125	129.0711		7.3526	B	131.58	92.5		IF
Potassium									NR
Selenium	75-125	2.6579		0.6316	B	2.63	77.0		IF
Silver	75-125	10.2974		0.5263	U	13.16	78.2		IF
Sodium									NR
Thallium	75-125	12.2105		0.5263	U	13.16	92.8		IF
Vanadium	75-125	140.6500		19.5842		131.58	92.0		IF
Zinc	75-125	289.7684		168.5632		131.58	92.1		IF
Cyanide	75-125	6.3737		0.6579	U	6.58	96.9		AS

Comments:



17744  
SDGMCJE !!

$$U = \langle IDL \rangle$$

## Graphite Furnace Spike Recovery Evaluation Form

Sample ID	Instr. Level Result $\mu\text{g/L}$	PDS Recovery %	Diluted Result $\mu\text{g/L}$	Diluted PDS Recovery %	MSA Result (if needed) $\mu\text{g/L}$	Final Result Reported $\text{mg/kg}$	
MCJE11	301 (hi)	53 (hi)	5.7x50	90 hi CV		113.4	100-230-
12	154	62	12.4x10	84	7.9x20	62.7	95.0
13	113.	71	24.8x10	74	5.0x20	33.3	72.0
14	198.2	70	8.7x20	79	5.7x40	105.5	145.0
15	870	-	18.7x100	74	7.3x200	349.	186.0
16	hi	hi	83.x100	125 (hi)		4810	4,200
17	107. ↓	50	8.6x10	84	3.9x20	35.7	30.0
18	90.5	84	6.9x10	89		18.2	32.0
18D	92.6	71	7.3x10	113		19.2	34.0
18S	104.1 (hi)	-	9.1x10			110.7	(90.0)
19	70.3	99				16.9	24.0
20	902 (hi)	92 (hi)	8.2x100	106		253.	413.0
21	(hi) ↓	hi	6.7x200	102		358.	360.0
22	87.4	101	8.3x10	116 (no rec)		22.2	66.0
23	105.4	100 ✓	7.8x10	113		20.2	40.0
26	60.6	87				17.8	44.0
MCJE11AL			5.8x50	115		113	
MCJE16AL			52.5x200	88		4810	

14P 1DL = 25.49 J

$$-U = < IDL$$

17744  
SDG MUSE II

ORIGINAL  
12/25

[illegible]

$$u = \langle IDL$$

17744  
SOGMOJE II

## Spike Recovery Evaluation Form

[illegible]

$$-u = <IDL$$

17744  
SDG-MJZ 11

[illegible]

8  
STANDARD ADDITION RESULTS

b Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Concentration Units: ug/L

EPA Sample No.	An	0 ADD ABS	1 ADD CON ABS	2 ADD CON ABS	3 ADD CON ABS	Final Conc.	r	Q
MCJE14	AS	0.043	5.00 0.077	10.00 0.095	15.00 0.123	8.9	0.9936	+
MCJE14	AS	0.043	5.00 0.072	10.00 0.100	15.00 0.123	8.3	0.9986	✓
MCJE15	AS	0.080	10.00 0.105	20.00 0.135	30.00 0.156	31.1	0.9979	+
MCJE16	AS	0.048	10.00 0.072	20.00 0.100	30.00 0.080?	45.5	0.7442	+
MCJE16	AS	0.049	10.00 0.076	20.00 0.099	30.00 0.076	57.1	0.6567	+
MCJE17	AS	0.040	10.00 0.067	20.00 0.094	30.00 0.116	16.1	0.9988	✓
MCJE18D	AS	0.060	15.00 0.095	30.00 0.118	45.00 0.160	27.8	0.9943	+
MCJE18D	AS	0.055	15.00 0.098	30.00 0.120	45.00 0.151	28.8	0.9917	+
MCJE21	AS	0.053	10.00 0.082	20.00 0.105	30.00 0.120	25.2	0.9903	+
MCJE21	AS	0.055	10.00 0.075	20.00 0.098	30.00 0.120	24.9	0.9996	✓
MCJE12	PB	0.022	5.00 0.036	10.00 0.049	15.00 0.064	7.9	0.9996	+
MCJE13	PB	0.015	5.00 0.032	10.00 0.046	15.00 0.062	5.0	0.9994	+
MCJE14	PB	0.016	5.00 0.028	10.00 0.044	15.00 0.056	5.7	0.9983	+
MCJE15	PB	0.020	5.00 0.036	10.00 0.050	15.00 0.063	7.3	0.9989	+
MCJE17	PB	0.012	5.00 0.027	10.00 0.043	15.00 0.058	3.9	0.9999	+
MCJE12	SE	0.003	5.00 0.019	10.00 0.025	15.00 0.041	1.7	0.9864	+
MCJE12	SE	0.005	5.00 0.007	10.00 0.028	15.00 0.035	0.9	0.9541	+
MCJE14	SE	0.005	5.00 0.016	10.00 0.029	15.00 0.044	1.5	0.9976	+
MCJE14	SE	0.012	5.00 0.024	10.00 0.042	15.00 0.055	3.8	0.9971	+
MCJE15	SE	0.004	5.00 0.015	10.00 0.030	15.00 0.040	1.5	0.9972	+

see  
KF's  
on  
pos  
for

## U.S. EPA - CLP

6  
DUPLICATES

EPA SAMPLE NO.

MCJE18D

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Level (low/med): LOW

% Solids for Sample: 76.0

% Solids for Duplicate: 73.4

Concentration Units (ug/L or mg/kg dry weight): MG/KG

Analyte	Control Limit	Sample (S)	C	Duplicate (D)	C	RPD	Q	M
Aluminum		3397.4634		3491.8000		2.7		P
Antimony		7.8347	U	7.8347	U			P
Arsenic	2.6	6.6579		7.3081		9.3		F
Barium	52.6	52.7026		51.8816	B	1.6		P
Beryllium		0.5105	B	0.5763	B	12.1		P
Cadmium		0.7895	U	0.7895	U			P
Calcium	1315.8	2108.6553		2130.3291		1.0		P
Chromium	2.6	10.6316		10.9921		3.3		P
Cobalt		6.3658	B	6.2395	B	2.0		P
Copper	6.6	9.9684		10.2184		2.5		P
Iron		10750.5322		10928.9844		1.6		P
Lead		181.5789		192.1053		5.6		F
Magnesium		936.6500	B	971.0816	B	3.6		P
Manganese		317.6579		310.3868		2.3		P
Mercury		0.1316	U	0.1316	U			CV
Nickel		7.3526	B	6.1237	B	18.2		P
Potassium		397.3369	B	384.2658	B	3.3		P
Selenium		0.6316	B	0.5263	U	200.0		F
Silver		0.5263	U	0.5263	U			P
Sodium		43.5237	B	49.1895	B	12.2		P
Thallium		0.5263	U	0.5263	U			F
Vanadium	13.2	19.5842		19.7158		0.7		P
Zinc		168.5632		169.0105		0.3		P
Cyanide		0.6579	U	0.6579	U			AS

REPORT

KEYSTONE D.C.# 17744-3- 13

U.S. EPA - CLP

9  
ICP SERIAL DILUTIONS

EPA SAMPLE NO.

MCJE18L

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Level (low/med): LOW

Concentration Units: ug/L

Analyte	Initial Sample Result (I)	C	Serial Dilution Result (S)	C	% Differ- ence	Q	M
Aluminum_	12910.36		12943.00		0.3		P
Antimony_	30.00	U	150.00	U			P
Arsenic_							NR
Barium_	200.27		200.55	B	0.1		P
Beryllium_	1.94	B	5.00	U	100.0		P
Cadmium_	3.00	U	15.00	U			P
Calcium_	3012.89		3317.10	B	3.8		P
Chromium_	40.40		39.20	B	3.0		P
Cobalt_	24.19	B	20.55	B	15.0		P *
Copper_	37.88		35.50	B	6.3		P
Iron_	40852.02		41144.55		0.7		P
Lead_							NR
Magnesium_	3559.27	B	3600.75	B	1.2		P
Manganese_	1207.10		1222.75		1.3		P
Mercury_							NR
Nickel_	27.94	B	110.00	U	100.0		P
Potassium_	1509.88	B	3610.00	U	100.0		P
Selenium_							NR
Silver_	2.00	U	10.00	U			P
Sodium_	165.39	B	240.90	B	45.7		P *
Thallium_							NR
Vanadium_	74.42		76.45	B	2.7		P
Zinc_	640.54		662.70		3.5		P

\*  $\left(\frac{S-I}{S}\right) < 10\%, \text{ no impact}$

## FURNACE-AA BENCH SHEET

Page \_\_\_\_ of \_\_\_\_

KEYSTONE LAB-HOUSTON

D.C. # \_\_\_\_\_ -08

ELEMENT Se  
DATE 1-23-92INSTRUMENT TSA-21  
ANALYST [signature]CRDL 5  
IDL 2Note: Use proper units and flags  
for final concentration

CUP NO	SAMPLE OR STD ID	CLIENT ID NO.	INITIAL DATA			DILUTION DATA				MSA DATA		DIG FACT	ANALYTE CONC	
			AVG CONC	RPD	ANAL SPIKE REC %	DIL FACT	AVG CONC	RPD	ANAL SPIKE REC %	DIL FACT	CALC CONC INTCP		UG/L	MG/KG
27		A	11.1	11.44	860%								—	—
28		MCSE20	3.1										—	8.6
29		A	9.3	3.01	62%								—	—
30		MCSE21	0.9										—	8.4
31		A	8.1	9.62	81%								—	—
32		MCSE22	2.7			lab misread as 2.7; should be 2.2							—	0.8
33		A	7.6	16.71	540%								—	—
34		MCSE23	1.9										—	0.4
35		A	8.4	14.76	86%								—	—
36	000132	CCV	47.8	4.28									47.8 50.0	960%
37		CCB	0.6										2	—
38		MCSE24	0.7			1X	2.8							0.6
39		A	14.2	40.8		12	12.3	20.08	85%				—	—



## Se-S As-S

SN= 000031

Conc 1 008.7 HIGH  
 Mean 008.7 HIGH  
 P/H 0.047 0.254  
 Conc 2 007.6 HIGH  
 Mean 008.1 HIGH  
 /H 0.054 0.237

14A

810% see  
dil

## Se-S As-S

Mean 008.1 HIGH  
 SD 000.8  
 RSD 09.62

## Se-S As-S

SN= 000032

Conc 1 001.0 040.2  
 Mean 001.0 040.2  
 P/H 0.006 0.133  
 Conc 2 003.5 037.5  
 Mean 002.2 038.8  
 P/H 0.005 0.133

15 MCSE 22

0.4W 7.8

## Se-S As-S

Mean 002.2 038.8  
 SD 001.8 001.9  
 RSD 80.00 04.92

## Se-S As-S

SN= 000033

Conc 1 006.7 060.4  
 Mean 006.7 060.4  
 P/H 0.052 0.189  
 Conc 2 008.5 057.5  
 Mean 007.6 058.9  
 P/H 0.052 0.198

15A

590% 1000%

## Se-S As-S

Mean 007.6 058.9  
 SD 001.3 002.1  
 RSD 16.71 03.48

## Se-S As-S

SN= 000034

Conc 1 002.4 021.1  
 Mean 002.4 021.1  
 P/H 0.005 0.075  
 Conc 2 001.5 021.9  
 Mean 001.9 021.5  
 P/H 0.003 0.088

16 MCSE 23

0.4 4.3

## Se-S As-S

Mean 001.9 021.5  
 SD 000.6 000.6  
 RSD 33.68 02.60

## Se-S As-S

SN= 000035

Conc 1 009.5 046.0  
 Mean 009.5 046.0  
 P/H 0.059 0.186  
 Conc 2 007.7 040.8  
 Mean 008.6 043.4  
 P/H 0.038 0.168

16A

860% 1100%

## Se-S As-S

Mean 008.6 043.4  
 SD 001.3 003.7  
 RSD 14.76 08.45

000144

Pb. P.E. S60

H92-01.302

2/26/92

T.N

14:15

-1.6  
-1.4  
-1.5AV *ccs* 1.0u  
-11.85CV

8.6  
9.6  
9.1AV *oil* (spk) *ox*  
8.21CV

110  $\frac{g}{g}$  MCJE 18 (spk)

8.2  
8.2

8.2AV *oil* 13 (100x)

(8.2)(100)(2) = 164 mg/kg

0.75CV

29.2

29.5

29.3AV +20

105.5  $\frac{g}{g}$

MCJE 20A

0.60CV

6.5

6.8

6.7AV *oil* 14 (200x)

(6.7)(200)(2) = 268 mg/kg

2.77CV

27.3

26.8

27.0AV +20

101.5  $\frac{g}{g}$

MCJE 21A

1.48CV

8.6

8.1

8.3AV *oil* 15 (10x)

(8.3)(10)(2) = 16.6 mg/kg

4.34CV

31.6

31.4

31.5AV +20

116  $\frac{g}{g}$  NO MSA performed

MCJE 22A

0.44CV

7.5

8.2

7.8AV *oil* 16 (10x)

(7.8)(10)(2) = 15.6 mg/kg

6.24CV

31.3

29.8

30.5AV +20

113.5  $\frac{g}{g}$

MCJE 23A

3.58CV

52.3

49.9

51.1AV *ccv* 9

51.1/50.0 = 102.2  $\frac{g}{g}$

3.33CV

-1.3

-1.4

-1.3AV *ccs*

1.0u

-0.83CV

5.6

5.9

5.8AV *oil* 2 (50x)

(5.8)(50)(2) = 58 mg/kg

4.19CV

28.8

28.8

28.8AV +20

115  $\frac{g}{g}$

MCJE 11A

0.10CV

52.3

52.7

52.5AV *oil* 7 (200x)

(52.5)(200)(2) = 2100 mg/kg

0.46CV

MCJE 16

000168

(PL) ...

H92-07.302

2/26/92

T.N

71.9

ORIGINAL

76

68.1  
70.0AV +20 87.5 ~~ST~~ MCJE 16A

3.83CV

49.8

49.7

77

49.7AV ~~ccv~~ 49.7/50.0 = 99.4%

0.10CV

-1.4

-0.9

78

-1.1AV ~~ccv~~ 1.0u

-25.50CV

20X

003(2x)

79 0.022+0 MCJE 12

80 0.036+5 MCJE 12

81 0.049+10 MCJE 12

82 0.064+15 MCJE 12

corr = 0.999

Slope = 0.00278

Inter = 0.0219

$$x' = (7.87)(20)(2) = 31.5 \text{ mg/kg}$$

004(2x) 20X

83 0.015+0 MCJE 13

84 0.032+5 MCJE 13

85 0.046+10 MCJE 13

86 0.062+15 MCJE 13

corr = 0.999

Slope = 0.0031

Inter = 0.0155

$$x' = (5)(20)(2) = 20.0 \text{ mg/kg}$$

005(2x) 40X

87 0.016+0 MCJE 14

88 0.028+5 MCJE 14

89 0.044+10 MCJE 14

90 0.056+15 MCJE 14

corr = 0.998

Slope = 0.00272

Inter = 0.0156

$$x' = (5.73)(40)(2) = 45.8 \text{ mg/kg}$$

006(2x) 200X

91 0.020+0 MCJE 15

92 0.036+5 MCJE 15

93 0.050+10 MCJE 15

94 0.063+15 MCJE 15

corr = 0.998

Slope = 0.00286

Inter = 0.0208

$$x' = (7.27)(200)(2) = 290.0 \text{ mg/kg}$$

008(2x) 20X

95 0.012+0 MCJE 17

96 0.027+5 MCJE 17

97 0.043+10 MCJE 17

98 0.058+15 MCJE 17

corr = 0.999

Slope = 0.00308

Inter = 0.0119

$$x' = (3.86)(20)(2) = 15.4 \text{ mg/kg}$$

17:25

99 50.8 ~~ccv~~ 50.8/50.0 = 101.6%

100 -1.1 ~~ccv~~ 17:30 1.0u

000169

PE560

T. N. Lang

2/26/92

8:00

Daily Log Run For Pb by Furnace

PE

3.0

50.0

100.0

ICV(2x)

ICB

CCV

CCB

CRA

PBS

PBSA

LCSS(2x)

LCSSA

CCV

CCB

CCV 9:15

CCB 9:20

H92-01.30Z

002

+20

003

+20

004

+20

005

+20

006

+20

007

+20

008

+20

009

+20

010

+20

011

+20

012

+20

013

+20

014

+20

015

+20

016

+20

017

+20

018

+20

019

+20

33 +20

34 016

35 +20

36 CCV

37 CCB

38 017

39 +20

40 002(50x)

41 +20

42 003(10x)

43 +20

44 004(10x)

45 +20

46 005(20x)

47 +20

48 CCV

49 CCB

50 006(100x)

51 +20

52 007(100x)

53 +20

54 008(10x)

55 +20

56 009(10x)

57 +20

58 010(10x)

59 +20

60 CCV

61 CCB

62 011(10x)

013 (100x)

+20

6. min

65 014(20x)

66 +20

67 015(10x)

68 +20

69 016(10x)

70 +20

71 CCV

72 CCB

73 002(50x)

74 +20

75 007(20x)

76 +20

77 CCV

78 CCB

79 003(10x)

80 +5

81 +10

82 004(10x)

83 +5

84 +10

85 005(2x)

86 +5

87 +10

88 006(2x)

89 +5

90 +10

91 007(2x)

92 +5

93 +10

94 008(2x)

95 +5

96 +10

97 009(2x)

98 +5

99 +10

100 +15

17:35

CCV

17:30

CCB

dilution  
factors  
for  
MSA  
analysis

20x

20x

40x

200x

20x

000 70

REVIEWED BY:

URB EPA - CLP

(Red)

10

## Instrument Detection Limits (Quarterly)

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

ICP ID Number:

Date: 02/18/92

Flame AA ID Number:

Furnace AA ID Number: PE3030Z

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum			200		NR
Antimony			60		NR
Arsenic			10		NR
Barium			200		NR
Beryllium			5		NR
Cadmium			5		NR
Calcium			5000		NR
Chromium			10		NR
Cobalt			50		NR
Copper			25		NR
Iron			100		NR
Lead			3		NR
Magnesium			5000		NR
Manganese			15		NR
Mercury			0.2		NR
Nickel			40		NR
Potassium			5000		NR
Selenium	196.00	BZ	5	3.39E-01	NR
Silver			10		NR
Sodium			5000		NR
Thallium	276.80	BZ	10	2.0E-01	NR
Vanadium			50		NR
Zinc			20		NR

Comments:

FORM X - IN

000044

U.S. EPA - CLP

CPD(14)  
/nd

10

Instrument Detection Limits (Quarterly)

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

ICP ID Number:

TJA61

Date:

02/10/92

Flame AA ID Number:

Furnace AA ID Number:

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum_	308.20		200	14.0	P
Antimony_	206.80		60	30.0	P
Arsenic_			10		NR
Barium_	493.40		200	1.0	P
Beryllium_	313.00		5	1.0	P
Cadmium_	228.80		5	3.0	P
Calcium_	317.80		5000	7.0	P
Chromium_	267.70		10	3.0	P
Cobalt_	228.60		50	4.0	P
Copper_	324.70		25	3.0	P
Iron_	259.90		100	6.0	P
Lead_			3		NR
Magnesium_	279.00		5000	37.0	P
Manganese_	257.60		15	2.0	P
Mercury_			0.2		NR
Nickel_	231.60		40	22.0	P
Potassium_	766.40		5000	722.0	P
Selenium_			5		NR
Silver_	328.00		10	2.0	P
Sodium_	588.90		5000	30.0	P
Thallium_			10		NR
Vanadium_	292.40		50	4.0	P
Zinc_	213.80		20	2.0	P

Comments:

FORM X - IN

000045

## U.S. EPA - CLP

10

## Instrument Detection Limits (Quarterly)

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

ICP ID Number:

Date: 02/14/92

Flame AA ID Number:

Furnace AA ID Number: PE560

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum_			200		NR
Antimony_			60		NR
Arsenic_			10		NR
Barium_			200		NR
Beryllium_			5		NR
Cadmium_			5		NR
Calcium_			5000		NR
Chromium_			10		NR
Cobalt_			50		NR
Copper_			25		NR
Iron_			100		NR
Lead_	283.30	BD	3	1.0	F
Magnesium_			5000		NR
Manganese_			15		NR
Mercury_			0.2		NR
Nickel_			40		NR
Potassium_			5000		NR
Selenium_			5		NR
Silver_			10		NR
Sodium_			5000		NR
Thallium_			10		NR
Vanadium_			50		NR
Zinc_			20		NR

Comments:

FORM X - IN

000046

U.S. EPA - CLP

10

Instrument Detection Limits (Quarterly)

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

ICP ID Number:

Date: 02/05/92

Flame AA ID Number: MAS50A

Furnace AA ID Number:

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum			200		NR
Antimony			60		NR
Arsenic			10		NR
Barium			200		NR
Beryllium			5		NR
Cadmium			5		NR
Calcium			5000		NR
Chromium			10		NR
Cobalt			50		NR
Copper			25		NR
Iron			100		NR
Lead			3		NR
Magnesium			5000		NR
Manganese			15		NR
Mercury	253.70		0.2	0.2	CV
Nickel			40		NR
Potassium			5000		NR
Selenium			5		NR
Silver			10		NR
Sodium			5000		NR
Thallium			10		NR
Vanadium			50		NR
Zinc			20		NR

Comments:

FORM X - IN

000047



U.S. EPA - CLP

10

## Instrument Detection Limits (Quarterly)

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

ICP ID Number:

Date: 02/12/92

Flame AA ID Number:

Furnace AA ID Number: TJASH21

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum_	197.30	BS	200	2.0	NR
Antimony_			60		NR
Arsenic_			10		F
Barium_			200		NR
Beryllium_			5		NR
Cadmium_			5		NR
Calcium_			5000		NR
Chromium_			10		NR
Cobalt_			50		NR
Copper_			25		NR
Iron_	196.00	BS	100	2.0	NR
Lead_			3		NR
Magnesium_			5000		NR
Manganese_			15		NR
Mercury_			0.2		NR
Nickel_			40		NR
Potassium_			5000		NR
Selenium_			5		F
Silver_			10		NR
Sodium_			5000		NR
Thallium_			10		NR
Vanadium_			50		NR
Zinc_			20		NR

Comments:

FORM X - IN

000048

TPO: ☐ ACTION ☒ FYIRegion 3**INORGANIC REGIONAL DATA ASSESSMENT SUMMARY**

CASE NO. 17744 LABORATORY KEYTX  
SDG NO. MCPE02 (AQ) DATA USER HNU-S  
SOW 3/90 REVIEW COMPLETION DATE 5/5/92  
NO. OF SAMPLES 10 MATRIX AA  
REVIEWER ☐ ESD ☐ ESAT ☒ OTHER, CONTRACT/CONTRACTOR HNU-S

	ICP	AA	Hg	CYANIDE
1. HOLDING TIMES	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
2. INITIAL CALIBRATIONS	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
3. CONTINUING CALIBRATIONS	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
4. FIELD BLANKS ("F" = not applicable)	<u>X</u>	<u>0</u>	<u>0</u>	<u>0</u>
5. LABORATORY BLANKS	<u>A</u>	<u>X</u>	<u>0</u>	<u>0</u>
6. ICS	<u>0</u>			
7. LCS	<u>0</u>			
8. DUPLICATE ANALYSIS	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
9. MATRIX SPIKE	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
10. MSA		<u>X</u>		
11. SERIAL DILUTION	<u>X</u>			
12. SAMPLE VERIFICATION	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
13. REGIONAL QC ("F" = not applicable)	<u>M</u>	<u>X</u>	<u>0</u>	<u>0</u>
14. OVERALL ASSESSMENT	<u>X</u>	<u>X</u>	<u>0</u>	<u>0</u>

O = No problems or minor problems that do not affect data usability.

X = No more than about 5% of the data points are qualified as either estimated or unusable.

M = More than about 5% of the data points are qualified as estimated.

Z = More than about 5% of the data points are qualified as unusable.

A = DPO action requested; use in conjunction with one of the above codes.

DPO ACTION ITEMS: Aqueous MSA results for Se in 3 porewell samples  
were erratic, suggesting analytical malfunction or mispiking. Also, one Se result  
misreported. no significant / contractual problems

AREAS OF CONCERN: \_\_\_\_\_

TPO: ☒ ACTION ☒ FYIRegion 3**INORGANIC REGIONAL DATA ASSESSMENT SUMMARY**

CASE NO. 17744 LABORATORY Key TX  
 SDG NO. MCJE II DATA USER HMS  
 SOW 3/90 REVIEW COMPLETION DATE 5/5/92  
 NO. OF SAMPLES 14 MATRIX solid  
 REVIEWER ☐ ESD ☐ ESAT ☒ OTHER, CONTRACT/CONTRACTOR HMS

	ICP	AA	Hg	CYANIDE
1. HOLDING TIMES	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
2. INITIAL CALIBRATIONS	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
3. CONTINUING CALIBRATIONS	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
4. FIELD BLANKS ("F" = not applicable)	<u>X</u>	<u>X</u>	<u>0</u>	<u>0</u>
5. LABORATORY BLANKS	<u>X</u>	<u>X</u>	<u>0</u>	<u>0</u>
6. ICS	<u>0</u>			
7. LCS	<u>0</u>			
8. DUPLICATE ANALYSIS	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
9. MATRIX SPIKE	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
10. MSA		<u>X</u>		
11. SERIAL DILUTION	<u>0</u>			
12. SAMPLE VERIFICATION	<u>0</u>	<u>0</u>	<u>0</u>	<u>0</u>
13. REGIONAL QC ("F" = not applicable)	<u>X</u>	<u>0</u>	<u>0</u>	<u>0</u>
14. OVERALL ASSESSMENT	<u>X</u>	<u>X</u>	<u>0</u>	<u>0</u>

O = No problems or minor problems that do not affect data usability.

X = No more than about 5% of the data points are qualified as either estimated or unusable.

M = More than about 5% of the data points are qualified as estimated.

Z = More than about 5% of the data points are qualified as unusable.

A = DPO action requested; use in conjunction with one of the above codes.

DPO ACTION ITEMS:

lab miscalculated 6 Pb results and  
one Se result. An MSA analysis  
should have been performed for Pb in  
one sample where the PDS rec'd > 115 %.

AREAS OF CONCERN:

## FURNACE-AA LENCH SHEET

Page \_\_\_\_ of \_\_\_\_

KEYSTONE LAB-HOUSTON

D.C. # \_\_\_\_\_ -08

ELEMENT SeINSTRUMENT TSA-21CRDL 5

Note: Use proper units and flags

DATE 1-23-92ANALYST [Signature]IDL 2

for final concentration

CUP NO	SAMPLE OR STD ID	CLIENT ID NO.	INITIAL DATA			DILUTION DATA				MSA DATA		DIG FACT	ANALYTE CONC	
			AVG CONC	RPD	ANAL SPIKE REC %	DIL FACT	AVG CONC	RPD	ANAL SPIKE REC %	DIL FACT	CALC CONC INTCP		UG/L	MG/KG
27		A	11.1	11.44	860%								—	—
28		MCSE20	3.1										—	8.6
29		A	9.3	3.01	62%								—	—
30		MCSE21	0.9										—	0.4
31		A	8.2	9.62	81%								—	—
32		MCSE22	2.2										—	0.9
33		A	7.6	16.71	540%								—	—
34		MCSE23	1.9										—	0.4
35		A	8.4	14.76	860%								—	—
36	000132	CCW	47.8	4.28									47.8 50.5	960%
37		CCB	0.6										2	—
38		MCSE24	0.7			1K	2.8						—	0.6
39		A	14.2	40.8		12	12.3	10.8	85%				—	—

Metal Pb  
IDL 1 ug/l  
U = < IDL

17744  
SDG MJE II

Graphite Furnace Spike Recovery Evaluation Form

Sample ID	Instr. Level Result ug/l	PDS Recovery %	Diluted Result ug/l	Diluted PDS Recovery %	MSA Result (if needed) ug/l	Final Result Reported mg/kg	UF
see below * MJE II	301 (hi)	53 (hi)	5.7 x 50	90 hi CV		113.4	236
12	154	62	12.4 x 10	84	7.9 x 20	62.7	95
13	113	71	24.8 x 10	74	5.0 x 20	33.3	72
14	198.2	70	8.7 x 20	79	5.7 x 40	105.5	145
15	870	—	18.7 x 100	74	7.3 x 200	342	186
see below * 16	hi	hi	83. x 100	125 (hi)		4810	4,2
17	107. ↓	50	8.6 x 10	84	3.9 x 20	35.7	3
18	90.5	84	6.9 x 10	89	* 18.2	18.2	32
18D	92.6	71	7.3 x 10	113		19.2	34
18S	104.1 (hi)	—	9.1 x 10			110.7	(90
19	70.3	99				16.9	24
20	962 (hi)	92 (hi)	8.2 x 100	106		253	413
21	(hi) ↓	hi	6.7 x 200	102		358	360
22	87.4	101	8.3 x 10	116 (no RE)		22.2	66
23	105.4	100 ✓	7.8 x 10	113		20.2	40
26	60.6	87				17.8	64
MJE I AL			5.8 x 50	115		113	
MJE I 6 AL			52.5 x 200	88		4810	

circled results have been reported; lab did not take dilution factors into account when calculating results for MJE II, lab misplaced decimal.

1 CP IDL = 25 ug/l

Se-S As-S  
 SN= 000031  
 Conc 1 008.7 HIGH 14A  
 Mean 008.7 HIGH  
 P/H 0.047 0.254  
 Conc 2 007.6 HIGH  
 Mean 008.1 HIGH 810% see  
 /H 0.054 0.237 dil  
 Se-S As-S  
 Mean 008.1 HIGH  
 SD 000.8  
 RSD 09.62

Se-S As-S  
 SN= 000032  
 Conc 1 001.0 040.2  
 Mean 001.0 040.2  
 P/H 0.006 0.133  
 Conc 2 003.5 037.5  
 Mean 002.2 038.8  
 P/H 0.005 0.133  
 Se-S As-S  
 Mean 002.2 038.8  
 SD 001.8 001.9  
 RSD 80.00 04.92

Se-S As-S  
 SN= 000033  
 Conc 1 006.7 060.4  
 Mean 006.7 060.4  
 P/H 0.052 0.189  
 Conc 2 008.5 057.5  
 Mean 007.6 058.9  
 P/H 0.052 0.198  
 Se-S As-S  
 Mean 007.6 058.9  
 SD 001.3 002.1  
 RSD 16.71 03.48

Se-S As-S  
 SN= 000034  
 Conc 1 002.4 021.1  
 Mean 002.4 021.1  
 P/H 0.005 0.075  
 Conc 2 001.5 021.9  
 Mean 001.9 021.5  
 P/H 0.003 0.088  
 Se-S As-S  
 Mean 001.9 021.5  
 SD 000.6 000.6  
 RSD 33.68 02.60

Se-S As-S  
 SN= 000035  
 Conc 1 009.5 046.0  
 Mean 009.5 046.0  
 P/H 0.059 0.186  
 Conc 2 007.7 040.8  
 Mean 008.6 043.4  
 P/H 0.038 0.168  
 Se-S As-S  
 Mean 008.6 043.4  
 SD 001.3 003.7  
 RSD 14.76 08.45

15 MCSE 22

0.4W 7.8

15A

59% 100%

16 MCSE 23

0.4 4.3

16A

86% 1100%

000144

H92-07.302  
2/26/92  
T.N

71.9  
68.1  
76 70.0AV +20 87.5<sup>2</sup> MCJE 16A  
3.83CV  
49.8 15:35  
49.7  
77 49.7AV CCV<sup>9</sup> 49.7/50.0 = 99.4<sup>2</sup>  
0.10CV  
-1.4 15:40  
-0.9  
78 -1.1AVCCB 1.0u  
-25.50CV

mg/kg values (not  
dry-wt.  
corrected)

003(2x) 20X  
79 0.022+0 MCJE 12  
80 0.036+5 MCJE 12  
81 0.049+10 MCJE 12  
82 0.064+15 MCJE 12

corr = 0.999  
Slope = 0.00278  
Inter = 0.0219

$$x' = (7.87)(20)(2) = 31.5 \text{ mg/kg}$$

004(2x) 20X  
83 0.015+0 MCJE 13  
84 0.032+5 MCJE 13  
85 0.046+10 MCJE 13  
86 0.062+15 MCJE 13

corr = 0.999  
Slope = 0.0031  
Inter = 0.0155

$$x' = (5)(20)(2) = 20.0 \text{ mg/kg}$$

005(2x) 40X  
87 0.018+0 MCJE 14  
88 0.028+5 MCJE 14  
89 0.044+10 MCJE 14  
90 0.056+15 MCJE 14

corr = 0.998  
Slope = 0.00272  
Inter = 0.0156

$$x' = (5.73)(40)(2) = 45.8 \text{ mg/kg}$$

006(2x) 200X  
91 0.020+0 MCJE 15  
92 0.036+5 MCJE 15  
93 0.050+10 MCJE 15  
94 0.063+15 MCJE 15

corr = 0.998  
Slope = 0.00286  
Inter = 0.0208

$$x' = (7.27)(200)(2) = 290.0 \text{ mg/kg}$$

008(2x) 20X  
95 0.012+0 MCJE 17  
96 0.027+5 MCJE 17  
97 0.043+10 MCJE 17  
98 0.058+15 MCJE 17

corr = 0.999  
Slope = 0.00308  
Inter = 0.0119

$$x' = (3.86)(20)(2) = 15.4 \text{ mg/kg}$$

99 50.8 CCV<sup>10</sup> 50.8/50.0 = 101.6<sup>2</sup>  
100 -1.1 17:30  
1.0u

000169

PE560

T. N. Lang

Daily Log Run For CPB by Furnale

P

2/26/92 Bk 8:00

17:35

3.0

50.0

100.0 8:20

ICV (2x)

ICB 8:25

CCV 8:30

CCB 8:35

CRA

PBS

PRSA

LCSS (2x)

LCSSA

CCV 9:00

CCB 9:10

CCV 9:15

CCB 9:20

H92-01.302

002

+20

003

+20

004

+20

005

+20

006

+20

CCV 10:15

CCB 10:20

007

+20

008

+20

009

+20

010 (Dup)

+20

011 (Spk)

CCV 11:10

CCB 11:15

012

+20

013

+20

014

+20

015

+20

016

+20

017

+20

33 +20

34 016

35 +20

36 CCV 12:10

37 CCB 12:15

38 017

39 +20

40 002 (50x)

41 +20

42 003 (10x)

43 +20

44 004 (10x)

45 +20

46 005 (20x)

47 +20

48 CCV 13:10

49 CCB 13:15

50 006 (100x)

51 +20

52 (007 (100x))

53 +20

54 008 (10x)

55 +20

56 009 (10x)

57 +20

58 010 (10x)

59 +20

60 CCV 14:10

61 CCB 14:15

62 011 (Spk)

013 (100x)

+20

+5

+10

+15

65 014 (200x)

66 +20

67 015 (10x)

68 +20

69 016 (10x)

70 +20

71 CCV 15:05

72 CCB 15:10

73 002 (50x)

74 +20

75 007 (200x)

76 +20

77 CCV 15:35

78 CCB 15:40

79 003 (100x)

80 +20

81 +5

82 +10

83 +15

84 004 (10x)

85 +20

86 +5

87 +10

88 +15

89 005 (20x)

90 +20

91 +5

92 +10

93 +15

94 006 (100x)

95 +20

96 +5

97 +10

98 +15

99 008 (10x)

100 +20

101 +5

102 +10

103 +15

duplicate  
factor  
for  
MSA  
analysis

20X  
40X  
200X  
20X

000 20

REVIEWED BY:



Pb. P.E. S60

H92-01.302

2/26/92

T.N

14:15

61

-1.6  
-1.4  
-1.5AV *ccs* 1.0u  
-11.85CV

62

8.6  
9.6  
9.1AV *oil* (spk) (ox)  
8.21CV

110  $\frac{g}{g}$  MCJE 18 (spk)

63

8.2  
8.2  
8.2AV *oil* 13 (100x)  
0.75CV

(8.2)(100)(.2) = 164mg/kg  
MCJE 20

64

29.2  
29.5  
29.3AV +20  
0.60CV

105.5  $\frac{g}{g}$  MCJE 20A

65

6.5  
6.8  
6.7AV *oil* 14 (200x)  
2.77CV

(6.7)(200)(.2) = 268mg/kg  
MCJE 21

66

27.3  
26.8  
27.0AV +20  
1.48CV

101.5  $\frac{g}{g}$

MCJE 21A

67

8.6  
8.1  
8.3AV *oil* 15 (10x)  
4.34CV

(8.3)(10)(.2) = 16.6mg/kg  
MCJE 22

68

31.6  
31.4  
31.5AV +20  
0.44CV

116  $\frac{g}{g}$  NO MSA performed

MCJE 22A

69

7.5  
8.2  
7.8AV *oil* 16 (10x)  
6.24CV

(7.8)(10)(.2) = 15.6mg/kg  
MCJE 23

70

31.3  
29.8  
30.5AV +20  
3.58CV

113.5  $\frac{g}{g}$

MCJE 23A

71

52.3  
49.9  
51.1AV *ccv* 9  
3.33CV

51.1/50.0 = 102.2  $\frac{g}{g}$

72

-1.3  
-1.4  
-1.3AV *ccs*  
-0.83CV

1.0u

73

5.6  
5.9  
5.8AV *oil* 2 (50x)  
4.19CV

(5.8)(50)(.2) = 58mg/kg  
MCJE 11

74

28.8  
28.8  
28.8AV +20  
0.10CV

115  $\frac{g}{g}$

MCJE 11A

75

52.3  
52.7  
52.5AV *oil* 7 (200x)  
0.46CV

(52.5)(200)(.2) = 2100mg/kg  
MCJE 16  
000168

## APPENDIX B

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO. 17744

CHY02

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HO SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477446

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077446B51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/02/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	24	B
67-64-1-----	Acetone	14	
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	1	J

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

039

ORIG:

Re:

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY02

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477446

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077446B51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.      Date Analyzed: 02/02/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 0 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY02

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477446

Sample wt/vol: 1000 (g/mL) ML Lab File ID: G2J77446C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/03/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/06/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO. COMPOUND Q

108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl) Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy) Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

271

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY02

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477446

Sample wt/vol: 1000 (g/mL) ML Lab File ID: G2J77446C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/03/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/06/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl) Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy) Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

272

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY02

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477446  
Sample wt/vol: 1000 (g/mL) ML Lab File ID: G2J77446C21  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture:        decanted: (Y/N)        Date Extracted: 02/03/92  
Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/06/92  
Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
PC Cleanup: (Y/N) N pH:       

Number TICs found: 7 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	12.19	5	J
2.	LABORATORY ARTIFACT	12.27	11	BJ
3.	UNKNOWN HYDROCARBON	13.20	2	J
4.	UNKNOWN HYDROCARBON	13.70	3	J
5.	UNKNOWN HYDROCARBON	14.17	3	J
6.	UNKNOWN	14.65	3	J
7.	UNKNOWN	16.09	10	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY02

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477446

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 01/30/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 02/04/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U



1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY03

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477451

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077451B51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/02/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	37	B
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

050

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY03

Lab Name: COMPUCHEM.RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HO SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477451  
Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077451B51  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec.      Date Analyzed: 02/02/92  
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 1

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 76-13-1	ETHANE, 1,1,2-TRICHLORO-1,2,	4.82	10	JN

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

ORIGINAL  
Filed  
EPA SAMPLE NO.

CHY03

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477451

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077451B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
108-95-2	Phenol	10	U
111-44-4	bis(2-Chloroethyl) Ether	10	U
95-57-8	2-Chlorophenol	10	U
541-73-1	1,3-Dichlorobenzene	10	U
106-46-7	1,4-Dichlorobenzene	10	U
95-50-1	1,2-Dichlorobenzene	10	U
95-48-7	2-Methylphenol	10	U
108-60-1	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5	4-Methylphenol	10	U
621-64-7	N-Nitroso-Di-n-Propylamine	10	U
67-72-1	Hexachloroethane	10	U
98-95-3	Nitrobenzene	10	U
78-59-1	Isophorone	10	U
88-75-5	2-Nitrophenol	10	U
105-67-9	2,4-Dimethylphenol	10	U
111-91-1	bis(2-Chloroethoxy) Methane	10	U
120-83-2	2,4-Dichlorophenol	10	U
120-82-1	1,2,4-Trichlorobenzene	10	U
91-20-3	Naphthalene	10	U
106-47-8	4-Chloroaniline	10	U
87-68-3	Hexachlorobutadiene	10	U
59-50-7	4-Chloro-3-Methylphenol	10	U
91-57-6	2-Methylnaphthalene	10	U
77-47-4	Hexachlorocyclopentadiene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U
95-95-4	2,4,5-Trichlorophenol	25	U
91-58-7	2-Chloronaphthalene	10	U
88-74-4	2-Nitroaniline	25	U
131-11-3	Dimethyl Phthalate	10	U
208-96-8	Acenaphthylene	10	U
606-20-2	2,6-Dinitrotoluene	10	U
99-09-2	3-Nitroaniline	25	U
83-32-9	Acenaphthene	10	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

293

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY03

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477451

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077451B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	10	U
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

294

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY03

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HO SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477451  
Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077451B52  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92  
Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92  
Injection Volume:       2.0 (uL) Dilution Factor:       1.0  
GPC Cleanup: (Y/N) N pH:       

Number TICs found:   0   CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY03

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579H0 SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477451

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 01/30/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 02/04/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
---------	----------	---	---

319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY04

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477457

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077457B51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/02/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	57	B
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY04

Lab Name: COMPUCHEM.RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477457  
Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077457B51  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec.      Date Analyzed: 02/02/92  
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 1

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 76-13-1	ETHANE, 1,1,2-TRICHLORO-1,2,	4.82	10	JN



1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY04

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477457

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077457B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

PC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO. COMPOUND Q

108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl) Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy) Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

306

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY04

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HO SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477457

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077457B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	10	U
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

(initials)  
Date

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY04

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477457

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077457B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

PC Cleanup: (Y/N) N pH:       

Number TICs found: 0 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

ID  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY04

Lab Name: COMPUCHEM.RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477457

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 01/30/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 02/04/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.0029	J
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY05

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477458

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077458B51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/02/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
74-87-3	Chloromethane	10	U
74-83-9	Bromomethane	10	U
75-01-4	Vinyl Chloride	10	U
75-00-3	Chloroethane	10	U
75-09-2	Methylene Chloride	16	B
67-64-1	Acetone	10	U
75-15-0	Carbon Disulfide	10	U
75-35-4	1,1-Dichloroethene	10	U
75-34-3	1,1-Dichloroethane	10	U
540-59-0	1,2-Dichloroethene (total)	10	U
67-66-3	Chloroform	10	U
107-06-2	1,2-Dichloroethane	10	U
78-93-3	2-Butanone	10	U
71-55-6	1,1,1-Trichloroethane	10	U
56-23-5	Carbon Tetrachloride	10	U
75-27-4	Bromodichloromethane	10	U
78-87-5	1,2-Dichloropropane	10	U
10061-01-5	cis-1,3-Dichloropropene	10	U
79-01-6	Trichloroethene	10	U
124-48-1	Dibromochloromethane	10	U
79-00-5	1,1,2-Trichloroethane	10	U
71-43-2	Benzene	10	U
10061-02-6	Trans-1,3-Dichloropropene	10	U
75-25-2	Bromoform	10	U
108-10-1	4-Methyl-2-Pentanone	10	U
591-78-6	2-Hexanone	10	U
127-18-4	Tetrachloroethene	10	U
79-34-5	1,1,2,2-Tetrachloroethane	10	U
108-88-3	Toluene	10	U
108-90-7	Chlorobenzene	10	U
100-41-4	Ethylbenzene	10	U
100-42-5	Styrene	10	U
1330-20-7	Xylene (total)	10	U

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY05

Lab Name: COMPUCHEM.RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477458  
Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077458B51  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec.      Date Analyzed: 02/02/92  
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 0 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY05

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477458

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077458B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO. COMPOUND

108-95-2-----Phenol	10	U
111-44-4-----bis(2-Chloroethyl) Ether	10	U
95-57-8-----2-Chlorophenol	10	U
541-73-1-----1,3-Dichlorobenzene	10	U
106-46-7-----1,4-Dichlorobenzene	10	U
95-50-1-----1,2-Dichlorobenzene	10	U
95-48-7-----2-Methylphenol	10	U
108-60-1-----2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----4-Methylphenol	10	U
621-64-7-----N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----Hexachloroethane	10	U
98-95-3-----Nitrobenzene	10	U
78-59-1-----Isophorone	10	U
88-75-5-----2-Nitrophenol	10	U
105-67-9-----2,4-Dimethylphenol	10	U
111-91-1-----bis(2-Chloroethoxy) Methane	10	U
120-83-2-----2,4-Dichlorophenol	10	U
120-82-1-----1,2,4-Trichlorobenzene	10	U
91-20-3-----Naphthalene	10	U
106-47-8-----4-Chloroaniline	10	U
87-68-3-----Hexachlorobutadiene	10	U
59-50-7-----4-Chloro-3-Methylphenol	10	U
91-57-6-----2-Methylnaphthalene	10	U
77-47-4-----Hexachlorocyclopentadiene	10	U
88-06-2-----2,4,6-Trichlorophenol	10	U
95-95-4-----2,4,5-Trichlorophenol	25	U
91-58-7-----2-Chloronaphthalene	10	U
88-74-4-----2-Nitroaniline	25	U
131-11-3-----Dimethyl Phthalate	10	U
208-96-8-----Acenaphthylene	10	U
606-20-2-----2,6-Dinitrotoluene	10	U
99-09-2-----3-Nitroaniline	25	U
83-32-9-----Acenaphthene	10	U

FORM I SV-1

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SAMPLE DATA PACKAGE

17744 CHY 02

319

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY05

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HO SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477458

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077458B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	10	U
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b) Fluoranthene	10	U
207-08-9-----	Benzo(k) Fluoranthene	10	U
50-32-8-----	Benzo(a) Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	10	U
53-70-3-----	Dibenz(a,h) Anthracene	10	U
191-24-2-----	Benzo(g,h,i) Perylene	10	U

(1) - Cannot be separated from Diphenylamine



1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY05

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477458  
Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077458B52  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92  
Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92  
Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
GPC Cleanup: (Y/N) N pH:       

Number TICs found: 1

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	9.65	3	J

EPA SAMPLE NO.

CHY 05

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET  
Contract: 68  
SAS No.: 657

Lab Name: COMPUchem RTP

Lab Code: COMPU  
(soil/v

Lab Code: COMPU  
Matrix: (soil/water) WATER  
wt/vol: 1000

Matrix: \_\_\_\_\_  
Sample wt/vol: \_\_\_\_\_  
Pre: \_\_\_\_\_

decanted: (Y/N)  
(conc)

Extraction: (SepF/Cont/Sonc)  
Extract Volume: 2.0 (uL)

Concentrated Extract  
Injection Volume: 2.0 (uL)

Injection Vol  
GPC Cleanup: (Y/N)N

pH:

**SEPF**

0000(uL)

BD10083  
79HQ SDG No.: CHY02  
Lab Sample ID: 477458  
ID:

Lab File ID:

Lab File ID: \_\_\_\_\_  
Date Received: 01/29/92  
Extracted: 01/30/92

Date Received: 01/30/92  
Date Extracted: 01/30/92  
Analyzed: 02/04/92

Date Analyzed: 1

Dilution Factor: -  
Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO.	COMPOUND	CONCENTRATION UNIT (ug/L or ug/Kg) UG/L
319-84-6	alpha-BHC	0.050
319-85-7	beta-BHC	0.050
319-86-8	delta-BHC	0.050
58-89-9	gamma-BHC (Lindane)	0.050
76-44-8	Heptachlor	0.050
309-00-2	Aldrin	0.050
1024-57-3	Heptachlor epoxide	0.050
959-98-8	Endosulfan I	0.050
60-57-1	Dieldrin	0.050
72-55-9	4,4'-DDE	0.050
72-20-8	Endrin	0.050
33213-65-9	Endosulfan II	0.050
72-54-8	4,4'-DDD	0.050
1031-07-8	Endosulfan sulfate	0.050
50-29-3	4,4'-DDT	0.050
72-43-5	Methoxychlor	0.050
53494-70-5	Endrin ketone	0.050
7421-93-4	Endrin aldehyde	0.050
5103-71-9	alpha-Chlordane	0.050
5103-74-2	gamma-Chlordane	0.050
8001-35-2	Toxaphene	0.050
12674-11-2	Aroclor-1016	0.050
1104-28-2	Aroclor-1221	0.050
1104-28-2	Aroclor-1232	0.050
1104-28-2	Aroclor-1242	0.050
1104-28-2	Aroclor-1248	0.050
1104-28-2	Aroclor-1254	0.050

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY06

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477459

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077459B51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/02/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	32	B
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

FORM I VOA

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SAMPLE DATA PACKAGE

17744 CHY 02

079

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY06

Lab Name: COMPUCHEM.RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477459  
Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077459B51  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec.      Date Analyzed: 02/02/92  
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 1 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	LABORATORY ARTIFACT	21.30	13	J

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY06

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477459

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077459B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl) Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy) Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

333

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY06

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477459

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077459B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	2	J
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b) Fluoranthene	10	U
207-08-9-----	Benzo(k) Fluoranthene	10	U
50-32-8-----	Benzo(a) Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	10	U
53-70-3-----	Dibenz(a,h) Anthracene	10	U
191-24-2-----	Benzo(g,h,i) Perylene	10	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

334

ORIGINAL  
1/10/92

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY06

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477459

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077459B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

Number TICs found: 1

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.77	4	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY06

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477459

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 01/30/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 02/04/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.0071	JP
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U



1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY07

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477460

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077460B51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/02/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L      Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	54	B
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY07

Lab Name: COMPUCHEM RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477460  
Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077460B51  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec.      Date Analyzed: 02/02/92  
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 0 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY07

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 17744

SAS No.: 6579HQ

SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477460

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH077460B52

Level: (low/med) LOW

Date Received: 01/29/92

% Moisture:        decanted: (Y/N)       

Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 02/03/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl) Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy) Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

348

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY07

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477460

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077460B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/03/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	1	J
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	10	U
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

349

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY07

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477460  
Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077460B52  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92  
Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/03/92  
Injection Volume:       2.0 (uL) Dilution Factor:       1.0  
GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

Number TICs found:   0  

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

ID  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY07

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477460

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 01/30/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 02/04/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

319-84-6	alpha-BHC	0.050	U
319-85-7	beta-BHC	0.050	U
319-86-8	delta-BHC	0.050	U
58-89-9	gamma-BHC (Lindane)	0.050	U
76-44-8	Heptachlor	0.050	U
309-00-2	Aldrin	0.050	U
1024-57-3	Heptachlor epoxide	0.050	U
959-98-8	Endosulfan I	0.050	U
60-57-1	Dieldrin	0.10	U
72-55-9	4,4'-DDE	0.10	U
72-20-8	Endrin	0.10	U
33213-65-9	Endosulfan II	0.10	U
72-54-8	4,4'-DDD	0.10	U
1031-07-8	Endosulfan sulfate	0.10	U
50-29-3	4,4'-DDT	0.10	U
72-43-5	Methoxychlor	0.50	U
53494-70-5	Endrin ketone	0.10	U
7421-93-4	Endrin aldehyde	0.10	U
5103-71-9	alpha-Chlordane	0.050	U
5103-74-2	gamma-Chlordane	0.050	U
8001-35-2	Toxaphene	5.0	U
12674-11-2	Aroclor-1016	1.0	U
11104-28-2	Aroclor-1221	2.0	U
11141-16-5	Aroclor-1232	1.0	U
53469-21-9	Aroclor-1242	1.0	U
12672-29-6	Aroclor-1248	1.0	U
11097-69-1	Aroclor-1254	1.0	U
11096-82-5	Aroclor-1260	1.0	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY08

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477464

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077464B51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/02/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	22	B
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

098

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY08

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477464  
Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077464B51  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec.      Date Analyzed: 02/02/92  
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====



1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY08

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477464

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077464B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

108-95-2-----Phenol	10	U
111-44-4-----bis(2-Chloroethyl) Ether	10	U
95-57-8-----2-Chlorophenol	10	U
541-73-1-----1,3-Dichlorobenzene	10	U
106-46-7-----1,4-Dichlorobenzene	10	U
95-50-1-----1,2-Dichlorobenzene	10	U
95-48-7-----2-Methylphenol	10	U
108-60-1-----2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----4-Methylphenol	10	U
621-64-7-----N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----Hexachloroethane	10	U
98-95-3-----Nitrobenzene	10	U
78-59-1-----Isophorone	10	U
88-75-5-----2-Nitrophenol	10	U
105-67-9-----2,4-Dimethylphenol	10	U
111-91-1-----bis(2-Chloroethoxy)Methane	10	U
120-83-2-----2,4-Dichlorophenol	10	U
120-82-1-----1,2,4-Trichlorobenzene	10	U
91-20-3-----Naphthalene	10	U
106-47-8-----4-Chloroaniline	10	U
87-68-3-----Hexachlorobutadiene	10	U
59-50-7-----4-Chloro-3-Methylphenol	10	U
91-57-6-----2-Methylnaphthalene	10	U
77-47-4-----Hexachlorocyclopentadiene	10	U
88-06-2-----2,4,6-Trichlorophenol	10	U
95-95-4-----2,4,5-Trichlorophenol	25	U
91-58-7-----2-Chloronaphthalene	10	U
88-74-4-----2-Nitroaniline	25	U
131-11-3-----Dimethyl Phthalate	10	U
208-96-8-----Acenaphthylene	10	U
606-20-2-----2,6-Dinitrotoluene	10	U
99-09-2-----3-Nitroaniline	25	U
83-32-9-----Acenaphthene	10	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

362

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

CHY08

Lab Code: COMPU

Case No.: 17744

SAS No.: 6579HQ

SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477464

Sample wt/vol: 1000 (g/mL) ML

Lab File ID: GH077464B52

Level: (low/med) LOW

Date Received: 01/29/92

% Moisture:        decanted: (Y/N)       

Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL)

Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) UG/L

Q

51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	10	U
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b)Fluoranthene	10	U
207-08-9-----	Benzo(k)Fluoranthene	10	U
50-32-8-----	Benzo(a)Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----	Dibenz(a,h)Anthracene	10	U
191-24-2-----	Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY08

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477464  
Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077464B52  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92  
Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92  
Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
GPC Cleanup: (Y/N) N pH:       

Number TICs found: 0 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

ORIGINAL  
FBI

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY08

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477464

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 01/30/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 02/04/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.0046	JP
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.014	JP
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY09

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477465

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: C2R77465B51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/06/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	73	B
67-64-1-----	Acetone	34	
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

107

ORIGINAL  
(Red)

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY09

Lab Name: COMPUCHEM RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477465  
Sample wt/vol: 5.0 (g/mL) ML Lab File ID: C2R77465B51  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec.      Date Analyzed: 02/06/92  
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY09

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477465

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077465B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl) Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy) Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

375

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY09

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477465

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077465B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

## CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) UG/L

Q

51-28-5-----2,4-Dinitrophenol	25	U
100-02-7-----4-Nitrophenol	25	U
132-64-9-----Dibenzofuran	10	U
121-14-2-----2,4-Dinitrotoluene	10	U
84-66-2-----Diethylphthalate	10	U
7005-72-3-----4-Chlorophenyl-phenylether	10	U
86-73-7-----Fluorene	10	U
100-01-6-----4-Nitroaniline	25	U
534-52-1-----4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----N-Nitrosodiphenylamine (1)	10	U
101-55-3-----4-Bromophenyl-phenylether	10	U
118-74-1-----Hexachlorobenzene	10	U
87-86-5-----Pentachlorophenol	25	U
85-01-8-----Phenanthrene	10	U
120-12-7-----Anthracene	10	U
86-74-8-----Carbazole	10	U
84-74-2-----Di-n-Butylphthalate	10	U
206-44-0-----Fluoranthene	10	U
129-00-0-----Pyrene	10	U
85-68-7-----Butylbenzylphthalate	10	U
91-94-1-----3,3'-Dichlorobenzidine	10	U
56-55-3-----Benzo(a)Anthracene	10	U
218-01-9-----Chrysene	10	U
117-81-7-----bis(2-Ethylhexyl)Phthalate	10	U
117-84-0-----Di-n-Octyl Phthalate	10	U
205-99-2-----Benzo(b)Fluoranthene	10	U
207-08-9-----Benzo(k)Fluoranthene	10	U
50-32-8-----Benzo(a)Pyrene	10	U
193-39-5-----Indeno(1,2,3-cd)Pyrene	10	U
53-70-3-----Dibenz(a,h)Anthracene	10	U
191-24-2-----Benzo(g,h,i)Perylene	10	U

(1) - Cannot be separated from Diphenylamine



1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY09

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HO SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477465

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077465B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

Number TICs found: 3

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	TETRACHLOROETHANE	4.63	3	J
2.	UNKNOWN	4.77	4	J
3. 629-20-9	1,3,5,7-CYCLOOCTATETRAENE	4.28	47	J

381016A

(P00)

1D

PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY09

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579H0 SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477465

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 01/30/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 02/04/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.0050	JP
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY10

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477466

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077466B51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/03/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	7	BJ
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY10

Lab Name: COMPUCHEM.RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477466  
Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CN077466B51  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec.      Date Analyzed: 02/03/92  
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY10

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477466

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077466B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl) Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy) Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

391

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY10

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
 Matrix: (soil/water) WATER Lab Sample ID: 477466  
 Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077466B52  
 Level: (low/med) LOW Date Received: 01/29/92  
 % Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92  
 Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92  
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
 GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	10	U
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b) Fluoranthene	10	U
207-08-9-----	Benzo(k) Fluoranthene	10	U
50-32-8-----	Benzo(a) Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	10	U
53-70-3-----	Dibenz(a,h) Anthracene	10	U
191-24-2-----	Benzo(g,h,i) Perylene	10	U

(1) - Cannot be separated from Diphenylamine

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY10

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477466

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077466B52

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 02/01/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

Number TICs found: 2 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.73	6	J
2.	UNKNOWN	4.75	11	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY10

Lab Name: COMPUCHEM RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477466

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 01/30/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 02/04/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
319-84-6	alpha-BHC	0.050	U
319-85-7	beta-BHC	0.050	U
319-86-8	delta-BHC	0.050	U
58-89-9	gamma-BHC (Lindane)	0.050	U
76-44-8	Heptachlor	0.050	U
309-00-2	Aldrin	0.050	U
1024-57-3	Heptachlor epoxide	0.050	U
959-98-8	Endosulfan I	0.050	U
60-57-1	Dieldrin	0.10	U
72-55-9	4,4'-DDE	0.10	U
72-20-8	Endrin	0.10	U
33213-65-9	Endosulfan II	0.10	U
72-54-8	4,4'-DDD	0.0049	JP
1031-07-8	Endosulfan sulfate	0.10	U
50-29-3	4,4'-DDT	0.10	U
72-43-5	Methoxychlor	0.50	U
53494-70-5	Endrin ketone	0.10	U
7421-93-4	Endrin aldehyde	0.10	U
5103-71-9	alpha-Chlordane	0.050	U
5103-74-2	gamma-Chlordane	0.050	U
8001-35-2	Toxaphene	5.0	U
12674-11-2	Aroclor-1016	1.0	U
11104-28-2	Aroclor-1221	2.0	U
11141-16-5	Aroclor-1232	1.0	U
53469-21-9	Aroclor-1242	1.0	U
12672-29-6	Aroclor-1248	1.0	U
11097-69-1	Aroclor-1254	1.0	U
11096-82-5	Aroclor-1260	1.0	U



1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY11

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477472

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GR077472B54

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 29 Date Analyzed: 02/03/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:            (uL) Soil Aliquot Volume:            (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	14	U
74-83-9-----	Bromomethane	14	U
75-01-4-----	Vinyl Chloride	14	U
75-00-3-----	Chloroethane	14	U
75-09-2-----	Methylene Chloride	50	B
67-64-1-----	Acetone	54	B
75-15-0-----	Carbon Disulfide	14	U
75-35-4-----	1,1-Dichloroethene	14	U
75-34-3-----	1,1-Dichloroethane	14	U
540-59-0-----	1,2-Dichloroethene (total)	14	U
67-66-3-----	Chloroform	14	U
107-06-2-----	1,2-Dichloroethane	14	U
78-93-3-----	2-Butanone	14	U
71-55-6-----	1,1,1-Trichloroethane	3	J
56-23-5-----	Carbon Tetrachloride	14	U
75-27-4-----	Bromodichloromethane	14	U
78-87-5-----	1,2-Dichloropropane	14	U
10061-01-5-----	cis-1,3-Dichloropropene	14	U
79-01-6-----	Trichloroethene	14	U
124-48-1-----	Dibromochloromethane	14	U
79-00-5-----	1,1,2-Trichloroethane	14	U
71-43-2-----	Benzene	14	U
10061-02-6-----	Trans-1,3-Dichloropropene	14	U
75-25-2-----	Bromoform	14	U
108-10-1-----	4-Methyl-2-Pentanone	14	U
591-78-6-----	2-Hexanone	14	U
127-18-4-----	Tetrachloroethene	14	U
79-34-5-----	1,1,2,2-Tetrachloroethane	14	U
108-88-3-----	Toluene	14	U
108-90-7-----	Chlorobenzene	14	U
100-41-4-----	Ethylbenzene	14	U
100-42-5-----	Styrene	14	U
1330-20-7-----	Xylene (total)	14	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

048

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY11

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477472

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GDJ77472A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 29 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

GPC Cleanup: (Y/N) Y pH: 6.3

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND

108-95-2-----	Phenol	920	U
111-44-4-----	bis(2-Chloroethyl) Ether	920	U
95-57-8-----	2-Chlorophenol	920	U
541-73-1-----	1,3-Dichlorobenzene	920	U
106-46-7-----	1,4-Dichlorobenzene	920	U
95-50-1-----	1,2-Dichlorobenzene	920	U
95-48-7-----	2-Methylphenol	920	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	920	U
106-44-5-----	4-Methylphenol	920	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	920	U
67-72-1-----	Hexachloroethane	920	U
98-95-3-----	Nitrobenzene	920	U
78-59-1-----	Isophorone	920	U
88-75-5-----	2-Nitrophenol	920	U
105-67-9-----	2,4-Dimethylphenol	920	U
111-91-1-----	bis(2-Chloroethoxy) Methane	920	U
120-83-2-----	2,4-Dichlorophenol	920	U
120-82-1-----	1,2,4-Trichlorobenzene	920	U
91-20-3-----	Naphthalene	920	U
106-47-8-----	4-Chloroaniline	920	U
87-68-3-----	Hexachlorobutadiene	920	U
59-50-7-----	4-Chloro-3-Methylphenol	920	U
91-57-6-----	2-Methylnaphthalene	920	U
77-47-4-----	Hexachlorocyclopentadiene	920	U
88-06-2-----	2,4,6-Trichlorophenol	920	U
95-95-4-----	2,4,5-Trichlorophenol	2200	U
91-58-7-----	2-Chloronaphthalene	920	U
88-74-4-----	2-Nitroaniline	2200	U
131-11-3-----	Dimethyl Phthalate	920	U
208-96-8-----	Acenaphthylene	100	J
606-20-2-----	2,6-Dinitrotoluene	920	U
99-09-2-----	3-Nitroaniline	2200	U
83-32-9-----	Acenaphthene	180	J

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

435

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY11

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477472

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GDI77472A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 29 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

PC Cleanup: (Y/N) Y pH: 6.3

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

51-28-5-----	2,4-Dinitrophenol	2200	U
100-02-7-----	4-Nitrophenol	2200	U
132-64-9-----	Dibenzofuran	120	J
121-14-2-----	2,4-Dinitrotoluene	920	U
84-66-2-----	Diethylphthalate	180	J
7005-72-3-----	4-Chlorophenyl-phenylether	920	U
86-73-7-----	Fluorene	270	J
100-01-6-----	4-Nitroaniline	2200	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	2200	U
86-30-6-----	N-Nitrosodiphenylamine (1)	920	U
101-55-3-----	4-Bromophenyl-phenylether	920	U
118-74-1-----	Hexachlorobenzene	920	U
87-86-5-----	Pentachlorophenol	2200	U
85-01-8-----	Phenanthrene	2400	
120-12-7-----	Anthracene	550	J
86-74-8-----	Carbazole	250	J
84-74-2-----	Di-n-Butylphthalate	920	U
206-44-0-----	Fluoranthene	5600	
129-00-0-----	Pyrene	2900	
85-68-7-----	Butylbenzylphthalate	920	U
91-94-1-----	3,3'-Dichlorobenzidine	920	U
56-55-3-----	Benzo(a)Anthracene	2300	
218-01-9-----	Chrysene	1800	
117-81-7-----	bis(2-Ethylhexyl)Phthalate	280	J
117-84-0-----	Di-n-Octyl Phthalate	920	U
205-99-2-----	Benzo(b)Fluoranthene	3800	X
207-08-9-----	Benzo(k)Fluoranthene	3800	X
50-32-8-----	Benzo(a)Pyrene	1800	
193-39-5-----	Indeno(1,2,3-cd)Pyrene	1200	
53-70-3-----	Dibenz(a,h)Anthracene	410	J
191-24-2-----	Benzo(g,h,i)Perylene	1200	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

436

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY11

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477472

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GDJ77472A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 29 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

GPC Cleanup: (Y/N) Y pH: 6.3

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Number TICs found: 24

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.38	1100	J
2.	UNKNOWN	4.60	1300	J
3.	ALDOL	4.75	2100	ABJ
4.	UNKNOWN	5.20	1700	J
5.	UNKNOWN	5.83	1600	J
6.	BLANK CONTAMINANT	9.12	370	BJ
7.	BLANK CONTAMINANT	11.19	1100	BJ
8.	UNKNOWN SILOXANE	12.04	1100	J
9.	UNKNOWN CARBOXYLIC ACID	12.32	550	J
10.	UNKNOWN PAH	12.35	650	J
11.	UNKNOWN SILOXANE	12.82	1100	J
12.	UNKNOWN PAH	13.97	1300	J
13.	UNKNOWN PAH	14.07	650	J
14.	UNKNOWN PAH	14.12	1000	J
15.	UNKNOWN SILOXANE	14.19	2100	J
16.	LABORATORY ARTIFACT	14.52	1100	BJ
17.	UNKNOWN SILOXANE	14.80	2100	J
18.	UNKNOWN	15.05	740	J
19.	UNKNOWN SILOXANE	15.42	1600	J
20.	UNKNOWN PAH	15.75	920	J
21.	UNKNOWN	15.97	1300	J
22.	UNKNOWN SILOXANE	16.02	1300	J
23.	UNKNOWN	17.05	2100	J
24.	BENZOFLUORANTHENE	17.19	2300	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY11

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477472

Sample wt/vol: 30.40(g/ml)G

Lab File ID:

% Moisture: 29 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.3

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	Q
319-84-6-----	alpha-BHC	3.4
319-85-7-----	beta-BHC	2.4 U
319-86-8-----	delta-BHC	2.4 U
58-89-9-----	gamma-BHC (Lindane)	34 PE
76-44-8-----	Heptachlor	2.4 U
309-00-2-----	Aldrin	4.4 P
1024-57-3-----	Heptachlor epoxide	0.95 JP
959-98-8-----	Endosulfan I	2.4 U
60-57-1-----	Dieldrin	4.6 U
72-55-9-----	4,4'-DDE	4.6 U
72-20-8-----	Endrin	3.2 JP
33213-65-9-----	Endosulfan II	4.6 U
72-54-8-----	4,4'-DDD	2.8 JP
1031-07-8-----	Endosulfan sulfate	4.6 U
50-29-3-----	4,4'-DDT	4.6 U
72-43-5-----	Methoxychlor	7.8 JPB
53494-70-5-----	Endrin ketone	3.8 JP
7421-93-4-----	Endrin aldehyde	4.6 U
5103-71-9-----	alpha-Chlordane	5.3 P
5103-74-2-----	gamma-Chlordane	1.0 JP
8001-35-2-----	Toxaphene	240 U
12674-11-2-----	Aroclor-1016	46 U
11104-28-2-----	Aroclor-1221	93 U
11141-16-5-----	Aroclor-1232	46 U
53469-21-9-----	Aroclor-1242	46 U
12672-29-6-----	Aroclor-1248	46 U
11097-69-1-----	Aroclor-1254	46 U
11096-82-5-----	Aroclor-1260	46 U

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY11DL

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477472 D20

Sample wt/vol: 30.40(g/ml)G

Lab File ID:

% Moisture: 29 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/13/92

Injection Volume: 2.0(uL)

Dilution Factor: 2

GPC Cleanup: (Y/N) Y

pH: 6.3

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6-----	alpha-BHC	4.7	U
319-85-7-----	beta-BHC	4.7	U
319-86-8-----	delta-BHC	1.5	JPD
58-89-9-----	gamma-BHC (Lindane)	31	PD
76-44-8-----	Heptachlor	4.7	U
309-00-2-----	Aldrin	3.7	JPD
1024-57-3-----	Heptachlor epoxide	1.7	JPD
959-98-8-----	Endosulfan I	4.7	U
60-57-1-----	Dieldrin	9.2	U
72-55-9-----	4,4'-DDE	9.2	U
72-20-8-----	Endrin	3.9	JPD
33213-65-9-----	Endosulfan II	9.2	U
72-54-8-----	4,4'-DDD	1.8	JPD
1031-07-8-----	Endosulfan sulfate	9.2	U
50-29-3-----	4,4'-DDT	9.2	U
72-43-5-----	Methoxychlor	47	U
53494-70-5-----	Endrin ketone	9.2	U
7421-93-4-----	Endrin aldehyde	9.2	U
5103-71-9-----	alpha-Chlordane	4.4	JPD
5103-74-2-----	gamma-Chlordane	2.7	JD
8001-35-2-----	Toxaphene	470	U
12674-11-2-----	Aroclor-1016	92	U
11104-28-2-----	Aroclor-1221	190	U
11141-16-5-----	Aroclor-1232	92	U
53469-21-9-----	Aroclor-1242	92	U
12672-29-6-----	Aroclor-1248	92	U
11097-69-1-----	Aroclor-1254	92	U
11096-82-5-----	Aroclor-1260	92	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY12

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477476

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077476C18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 55 Date Analyzed: 01/31/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	22	U
74-83-9-----	Bromomethane	22	U
75-01-4-----	Vinyl Chloride	22	U
75-00-3-----	Chloroethane	22	U
75-09-2-----	Methylene Chloride	88	B
67-64-1-----	Acetone	96	B
75-15-0-----	Carbon Disulfide	22	U
75-35-4-----	1,1-Dichloroethene	22	U
75-34-3-----	1,1-Dichloroethane	22	U
540-59-0-----	1,2-Dichloroethene (total)	22	U
67-66-3-----	Chloroform	22	U
107-06-2-----	1,2-Dichloroethane	22	U
78-93-3-----	2-Butanone	22	U
71-55-6-----	1,1,1-Trichloroethane	22	U
56-23-5-----	Carbon Tetrachloride	22	U
75-27-4-----	Bromodichloromethane	22	U
78-87-5-----	1,2-Dichloropropane	22	U
10061-01-5-----	cis-1,3-Dichloropropene	22	U
79-01-6-----	Trichloroethene	22	U
124-48-1-----	Dibromochloromethane	22	U
79-00-5-----	1,1,2-Trichloroethane	22	U
71-43-2-----	Benzene	22	U
10061-02-6-----	Trans-1,3-Dichloropropene	22	U
75-25-2-----	Bromoform	22	U
108-10-1-----	4-Methyl-2-Pentanone	22	U
591-78-6-----	2-Hexanone	22	U
127-18-4-----	Tetrachloroethene	22	U
79-34-5-----	1,1,2,2-Tetrachloroethane	22	U
108-88-3-----	Toluene	22	U
108-90-7-----	Chlorobenzene	22	U
100-41-4-----	Ethylbenzene	22	U
100-42-5-----	Styrene	22	U
1330-20-7-----	Xylene (total)	22	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

060

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY12

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477476

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GR077476C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 55 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/10/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.4

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	Q
108-95-2	Phenol	75 J
111-44-4	bis(2-Chloroethyl) Ether	720 U
95-57-8	2-Chlorophenol	720 U
541-73-1	1,3-Dichlorobenzene	720 U
106-46-7	1,4-Dichlorobenzene	720 U
95-50-1	1,2-Dichlorobenzene	720 U
95-48-7	2-Methylphenol	720 U
108-60-1	2,2'-Oxybis(1-Chloropropane)	720 U
106-44-5	4-Methylphenol	95 J
621-64-7	N-Nitroso-Di-n-Propylamine	720 U
67-72-1	Hexachloroethane	720 U
98-95-3	Nitrobenzene	720 U
78-59-1	Isophorone	720 U
88-75-5	2-Nitrophenol	720 U
105-67-9	2,4-Dimethylphenol	720 U
111-91-1	bis(2-Chloroethoxy) Methane	720 U
120-83-2	2,4-Dichlorophenol	720 U
120-82-1	1,2,4-Trichlorobenzene	720 U
91-20-3	Naphthalene	720 U
106-47-8	4-Chloroaniline	720 U
87-68-3	Hexachlorobutadiene	720 U
59-50-7	4-Chloro-3-Methylphenol	720 U
91-57-6	2-Methylnaphthalene	720 U
77-47-4	Hexachlorocyclopentadiene	720 U
88-06-2	2,4,6-Trichlorophenol	720 U
95-95-4	2,4,5-Trichlorophenol	1700 U
91-58-7	2-Chloronaphthalene	720 U
88-74-4	2-Nitroaniline	1700 U
131-11-3	Dimethyl Phthalate	720 U
208-96-8	Acenaphthylene	720 U
606-20-2	2,6-Dinitrotoluene	720 U
99-09-2	3-Nitroaniline	1700 U
83-32-9	Acenaphthene	720 U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

513



1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY12

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477476

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GR077476C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 55 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/10/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

Cleanup: (Y/N) Y pH: 6.4

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

51-28-5-----	2,4-Dinitrophenol	1700	U
100-02-7-----	4-Nitrophenol	1700	U
132-64-9-----	Dibenzofuran	720	U
121-14-2-----	2,4-Dinitrotoluene	720	U
84-66-2-----	Diethylphthalate	720	U
7005-72-3-----	4-Chlorophenyl-phenylether	720	U
86-73-7-----	Fluorene	720	U
100-01-6-----	4-Nitroaniline	1700	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1700	U
86-30-6-----	N-Nitrosodiphenylamine (1)	720	U
101-55-3-----	4-Bromophenyl-phenylether	720	U
118-74-1-----	Hexachlorobenzene	720	U
87-86-5-----	Pentachlorophenol	1700	U
85-01-8-----	Phenanthrene	150	J
120-12-7-----	Anthracene	720	U
86-74-8-----	Carbazole	720	U
84-74-2-----	Di-n-Butylphthalate	79	J
206-44-0-----	Fluoranthene	360	J
129-00-0-----	Pyrene	320	J
85-68-7-----	Butylbenzylphthalate	76	J
91-94-1-----	3,3'-Dichlorobenzidine	720	U
56-55-3-----	Benzo(a)Anthracene	210	J
218-01-9-----	Chrysene	210	J
117-81-7-----	bis(2-Ethylhexyl)Phthalate	350	J
117-84-0-----	Di-n-Octyl Phthalate	720	U
205-99-2-----	Benzo(b)Fluoranthene	440	JX
207-08-9-----	Benzo(k)Fluoranthene	440	JX
50-32-8-----	Benzo(a)Pyrene	180	J
193-39-5-----	Indeno(1,2,3-cd)Pyrene	88	J
53-70-3-----	Dibenz(a,h)Anthracene	720	U
191-24-2-----	Benzo(g,h,i)Perylene	87	J

(1) - Cannot be separated from Diphenylamine

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY12

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477476

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GR077476C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 55 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/10/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.4

Number TICs found: 22

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.52	1300	J
2.	UNKNOWN	4.73	1200	J
3.	ALDOL	4.87	1100	ABJ
4.	ALDOL	5.00	1200	AJ
5.	UNKNOWN	5.32	870	J
6.	UNKNOWN	5.53	440	J
7.	UNKNOWN	5.97	1500	J
8.	UNKNOWN	7.60	510	J
9.	UNKNOWN	8.28	290	J
10. 70-55-3	BENZENESULFONAMIDE, 4-METHYL	10.80	360	JN
11.	UNKNOWN	12.35	360	J
12.	UNKNOWN	12.39	510	J
13. 57-10-3	HEXADECANOIC ACID	12.44	730	JN
14.	UNKNOWN	13.44	220	J
15.	UNKNOWN HYDROCARBON	14.22	290	J
16.	UNKNOWN HYDROCARBON	15.17	510	J
17.	UNKNOWN	15.64	220	J
18.	UNKNOWN	16.10	730	J
19.	UNKNOWN	16.55	730	J
20.	UNKNOWN	16.85	660	J
21.	UNKNOWN	17.22	3000	J
22.	UNKNOWN	18.85	2200	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY12

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477476

Sample wt/vol: 30.30(g/ml)G

Lab File ID:

% Moisture: 55 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.4

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6-----	alpha-BHC	0.38	JP
319-85-7-----	beta-BHC	3.7	U
319-86-8-----	delta-BHC	3.7	U
58-89-9-----	gamma-BHC (Lindane)	3.7	U
76-44-8-----	Heptachlor	0.13	JPB
309-00-2-----	Aldrin	1.1	JP
1024-57-3-----	Heptachlor epoxide	3.7	U
959-98-8-----	Endosulfan I	3.7	U
60-57-1-----	Dieldrin	7.3	U
72-55-9-----	4,4'-DDE	4.4	JP
72-20-8-----	Endrin	7.6	
33213-65-9-----	Endosulfan II	2.5	JP
72-54-8-----	4,4'-DDD	4.8	J
1031-07-8-----	Endosulfan sulfate	4.7	JP
50-29-3-----	4,4'-DDT	3.2	JP
72-43-5-----	Methoxychlor	15	JPB
53494-70-5-----	Endrin ketone	7.3	U
7421-93-4-----	Endrin aldehyde	2.6	JP
5103-71-9-----	alpha-Chlordane	4.1	P
5103-74-2-----	gamma-Chlordane	3.5	JP
8001-35-2-----	Toxaphene	370	U
12674-11-2-----	Aroclor-1016	73	U
11104-28-2-----	Aroclor-1221	150	U
11141-16-5-----	Aroclor-1232	73	U
53469-21-9-----	Aroclor-1242	73	U
12672-29-6-----	Aroclor-1248	73	U
11097-69-1-----	Aroclor-1254	73	U
11096-82-5-----	Aroclor-1260	73	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY13

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477483

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077483A18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 36 Date Analyzed: 01/31/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	Chloromethane	16	U
74-83-9	Bromomethane	16	U
75-01-4	Vinyl Chloride	16	U
75-00-3	Chloroethane	16	U
75-09-2	Methylene Chloride	67	B
67-64-1	Acetone	54	B
75-15-0	Carbon Disulfide	16	U
75-35-4	1,1-Dichloroethene	16	U
75-34-3	1,1-Dichloroethane	16	U
540-59-0	1,2-Dichloroethene (total)	16	U
67-66-3	Chloroform	16	U
107-06-2	1,2-Dichloroethane	16	U
78-93-3	2-Butanone	16	U
71-55-6	1,1,1-Trichloroethane	16	U
56-23-5	Carbon Tetrachloride	16	U
75-27-4	Bromodichloromethane	16	U
78-87-5	1,2-Dichloropropane	16	U
10061-01-5	cis-1,3-Dichloropropene	16	U
79-01-6	Trichloroethene	16	U
124-48-1	Dibromochloromethane	16	U
79-00-5	1,1,2-Trichloroethane	16	U
71-43-2	Benzene	16	U
10061-02-6	Trans-1,3-Dichloropropene	16	U
75-25-2	Bromoform	16	U
108-10-1	4-Methyl-2-Pentanone	16	U
591-78-6	2-Hexanone	16	U
127-18-4	Tetrachloroethene	16	U
79-34-5	1,1,2,2-Tetrachloroethane	16	U
108-88-3	Toluene	16	U
108-90-7	Chlorobenzene	16	U
100-41-4	Ethylbenzene	16	U
100-42-5	Styrene	16	U
1330-20-7	Xylene (total)	16	U

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY13

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477483

Sample wt/vol: 30.3 (g/mL) G Lab File ID: G2J77483A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 36 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

108-95-2-----Phenol	510	U
111-44-4-----bis(2-Chloroethyl) Ether	510	U
95-57-8-----2-Chlorophenol	510	U
541-73-1-----1,3-Dichlorobenzene	510	U
106-46-7-----1,4-Dichlorobenzene	510	U
95-50-1-----1,2-Dichlorobenzene	510	U
95-48-7-----2-Methylphenol	510	U
108-60-1-----2,2'-Oxybis(1-Chloropropane)	510	U
106-44-5-----4-Methylphenol	510	U
621-64-7-----N-Nitroso-Di-n-Propylamine	510	U
67-72-1-----Hexachloroethane	510	U
98-95-3-----Nitrobenzene	510	U
78-59-1-----Isophorone	510	U
88-75-5-----2-Nitrophenol	510	U
105-67-9-----2,4-Dimethylphenol	510	U
111-91-1-----bis(2-Chloroethoxy)Methane	510	U
120-83-2-----2,4-Dichlorophenol	510	U
120-82-1-----1,2,4-Trichlorobenzene	510	U
91-20-3-----Naphthalene	510	U
106-47-8-----4-Chloroaniline	510	U
87-68-3-----Hexachlorobutadiene	510	U
59-50-7-----4-Chloro-3-Methylphenol	510	U
91-57-6-----2-Methylnaphthalene	510	U
77-47-4-----Hexachlorocyclopentadiene	510	U
88-06-2-----2,4,6-Trichlorophenol	510	U
95-95-4-----2,4,5-Trichlorophenol	1200	U
91-58-7-----2-Chloronaphthalene	510	U
88-74-4-----2-Nitroaniline	1200	U
131-11-3-----Dimethyl Phthalate	510	U
208-96-8-----Acenaphthylene	510	U
606-20-2-----2,6-Dinitrotoluene	510	U
99-09-2-----3-Nitroaniline	1200	U
83-32-9-----Acenaphthene	510	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

579

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY13

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477483

Sample wt/vol: 30.3 (g/mL) G Lab File ID: G2J77483A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 36 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND Q

51-28-5-----	2,4-Dinitrophenol	1200	U
100-02-7-----	4-Nitrophenol	1200	U
132-64-9-----	Dibenzofuran	510	U
121-14-2-----	2,4-Dinitrotoluene	510	U
84-66-2-----	Diethylphthalate	76	J
7005-72-3-----	4-Chlorophenyl-phenylether	510	U
86-73-7-----	Fluorene	510	U
100-01-6-----	4-Nitroaniline	1200	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1200	U
86-30-6-----	N-Nitrosodiphenylamine (1)	510	U
101-55-3-----	4-Bromophenyl-phenylether	510	U
118-74-1-----	Hexachlorobenzene	510	U
87-86-5-----	Pentachlorophenol	1200	U
85-01-8-----	Phenanthrene	72	J
120-12-7-----	Anthracene	510	U
86-74-8-----	Carbazole	510	U
84-74-2-----	Di-n-Butylphthalate	510	U
206-44-0-----	Fluoranthene	150	J
129-00-0-----	Pyrene	130	J
85-68-7-----	Butylbenzylphthalate	510	U
91-94-1-----	3,3'-Dichlorobenzidine	510	U
56-55-3-----	Benzo(a) Anthracene	100	J
218-01-9-----	Chrysene	100	J
117-81-7-----	bis(2-Ethylhexyl) Phthalate	53	J
117-84-0-----	Di-n-Octyl Phthalate	510	U
205-99-2-----	Benzo(b) Fluoranthene	180	JX
207-08-9-----	Benzo(k) Fluoranthene	180	JX
50-32-8-----	Benzo(a) Pyrene	110	J
193-39-5-----	Indeno(1,2,3-cd) Pyrene	86	J
53-70-3-----	Dibenz(a,h) Anthracene	510	U
191-24-2-----	Benzo(g,h,i) Perylene	91	J

(1) - Cannot be separated from Diphenylamine

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY13

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477483

Sample wt/vol: 30.3 (g/mL) G Lab File ID: G2J77483A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 36 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.5

Number TICs found: 27

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.40	310	J
2.	UNKNOWN	4.62	1500	J
3.	ALDOL	4.77	1400	ABJ
4.	ALDOL	4.88	210	AJ
5.	UNKNOWN	5.87	1900	J
6.	BLANK CONTAMINANT	9.12	210	BJ
7.	UNKNOWN CARBOXYLIC ACID	9.84	210	J
8.	BLANK CONTAMINANT	10.24	360	BJ
9.	BLANK CONTAMINANT	11.20	570	BJ
10.	UNKNOWN SILOXANE	12.05	570	J
11.	UNKNOWN	12.59	150	J
12.	UNKNOWN SILOXANE	12.84	520	J
13.	BLANK CONTAMINANT	13.54	1000	BJ
14.	UNKNOWN SILOXANE	14.20	670	J
15.	LABORATORY ARTIFACT	14.54	210	BJ
16.	UNKNOWN SILOXANE	14.82	620	J
17.	UNKNOWN	15.07	310	J
18.	UNKNOWN SILOXANE	15.44	520	J
19.	UNKNOWN SILOXANE	16.04	820	J
20.	UNKNOWN	16.42	460	J
21.	UNKNOWN	16.47	260	J
22.	UNKNOWN	16.57	210	J
23.	UNKNOWN	16.72	930	J
24.	UNKNOWN	17.07	520	J
25.	UNKNOWN	17.15	770	J
26.	UNKNOWN	23.39	570	J
27.	UNKNOWN	23.45	310	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY13

Lab Name: COMPUCHEM RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477483

Sample wt/vol: 30.20(g/ml)G

Lab File ID:

% Moisture: 36 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.5

Sulfur Cleanup: (Y/N) N

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

319-84-6-----	alpha-BHC	2.6	U
319-85-7-----	beta-BHC	2.6	U
319-86-8-----	delta-BHC	2.6	U
58-89-9-----	gamma-BHC (Lindane)	2.6	U
76-44-8-----	Heptachlor	0.14	JPB
309-00-2-----	Aldrin	0.33	J
1024-57-3-----	Heptachlor epoxide	0.24	JP
959-98-8-----	Endosulfan I	2.6	U
60-57-1-----	Dieldrin	0.46	JP
72-55-9-----	4,4'-DDE	5.1	U
72-20-8-----	Endrin	1.6	JP
33213-65-9-----	Endosulfan II	0.32	JP
72-54-8-----	4,4'-DDD	5.1	U
1031-07-8-----	Endosulfan sulfate	5.1	U
50-29-3-----	4,4'-DDT	5.1	U
72-43-5-----	Methoxychlor	2.7	JPB
53494-70-5-----	Endrin ketone	0.65	JP
7421-93-4-----	Endrin aldehyde	5.1	U
5103-71-9-----	alpha-Chlordane	0.26	JP
5103-74-2-----	gamma-Chlordane	0.13	J
8001-35-2-----	Toxaphene	260	U
12674-11-2-----	Aroclor-1016	51	U
11104-28-2-----	Aroclor-1221	100	U
11141-16-5-----	Aroclor-1232	51	U
53469-21-9-----	Aroclor-1242	51	U
12672-29-6-----	Aroclor-1248	51	U
11097-69-1-----	Aroclor-1254	51	U
11096-82-5-----	Aroclor-1260	51	U



1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY14

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477484

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GR077484A18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 56 Date Analyzed: 02/05/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:            (uL) Soil Aliquot Volume:            (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	23	U
74-83-9-----	Bromomethane	23	U
75-01-4-----	Vinyl Chloride	23	U
75-00-3-----	Chloroethane	23	U
75-09-2-----	Methylene Chloride	54	B
67-64-1-----	Acetone	39	B
75-15-0-----	Carbon Disulfide	23	U
75-35-4-----	1,1-Dichloroethene	23	U
75-34-3-----	1,1-Dichloroethane	23	U
540-59-0-----	1,2-Dichloroethene (total)	23	U
67-66-3-----	Chloroform	23	U
107-06-2-----	1,2-Dichloroethane	23	U
78-93-3-----	2-Butanone	23	U
71-55-6-----	1,1,1-Trichloroethane	23	U
56-23-5-----	Carbon Tetrachloride	23	U
75-27-4-----	Bromodichloromethane	23	U
78-87-5-----	1,2-Dichloropropane	23	U
10061-01-5-----	cis-1,3-Dichloropropene	23	U
79-01-6-----	Trichloroethene	23	U
124-48-1-----	Dibromochloromethane	23	U
79-00-5-----	1,1,2-Trichloroethane	23	U
71-43-2-----	Benzene	23	U
10061-02-6-----	Trans-1,3-Dichloropropene	23	U
75-25-2-----	Bromoform	23	U
108-10-1-----	4-Methyl-2-Pentanone	23	U
591-78-6-----	2-Hexanone	23	U
127-18-4-----	Tetrachloroethene	23	U
79-34-5-----	1,1,2,2-Tetrachloroethane	23	U
108-88-3-----	Toluene	23	U
108-90-7-----	Chlorobenzene	23	U
100-41-4-----	Ethylbenzene	23	U
100-42-5-----	Styrene	23	U
1330-20-7-----	Xylene (total)	23	U

SAMPLE DATA PACKAGE

FORM I VOA  
17744 CHY 11

3/90

084

ORIGINAL  
Data

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY14

Lab Name: COMPUCHEM.RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
Matrix: (soil/water) SOIL Lab Sample ID: 477484  
Sample wt/vol: 30.4 (g/mL) G Lab File ID: GR077484C21  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: 56 decanted: (Y/N) N Date Extracted: 02/04/92  
Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/11/92  
Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
GPC Cleanup: (Y/N) Y pH: 6.2

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND		
108-95-2	Phenol	740	U
111-44-4	bis(2-Chloroethyl) Ether	740	U
95-57-8	2-Chlorophenol	740	U
541-73-1	1,3-Dichlorobenzene	740	U
106-46-7	1,4-Dichlorobenzene	740	U
95-50-1	1,2-Dichlorobenzene	740	U
95-48-7	2-Methylphenol	740	U
108-60-1	2,2'-Oxybis(1-Chloropropane)	740	U
106-44-5	4-Methylphenol	740	U
621-64-7	N-Nitroso-Di-n-Propylamine	740	U
67-72-1	Hexachloroethane	740	U
98-95-3	Nitrobenzene	740	U
78-59-1	Isophorone	740	U
88-75-5	2-Nitrophenol	740	U
105-67-9	2,4-Dimethylphenol	740	U
111-91-1	bis(2-Chloroethoxy) Methane	740	U
120-83-2	2,4-Dichlorophenol	740	U
120-82-1	1,2,4-Trichlorobenzene	740	U
91-20-3	Naphthalene	740	U
106-47-8	4-Chloroaniline	740	U
87-68-3	Hexachlorobutadiene	740	U
59-50-7	4-Chloro-3-Methylphenol	740	U
91-57-6	2-Methylnaphthalene	740	U
77-47-4	Hexachlorocyclopentadiene	740	U
88-06-2	2,4,6-Trichlorophenol	740	U
95-95-4	2,4,5-Trichlorophenol	1800	U
91-58-7	2-Chloronaphthalene	740	U
88-74-4	2-Nitroaniline	1800	U
131-11-3	Dimethyl Phthalate	740	U
208-96-8	Acenaphthylene	740	U
606-20-2	2,6-Dinitrotoluene	740	U
99-09-2	3-Nitroaniline	1800	U
83-32-9	Acenaphthene	740	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

645

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY14

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477484

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GR077484C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 56 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/11/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

IPC Cleanup: (Y/N) Y pH: 6.2

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND		Q
51-28-5-----	2,4-Dinitrophenol	1800	U
100-02-7-----	4-Nitrophenol	1800	U
132-64-9-----	Dibenzofuran	740	U
121-14-2-----	2,4-Dinitrotoluene	740	U
84-66-2-----	Diethylphthalate	740	U
7005-72-3-----	4-Chlorophenyl-phenylether	740	U
86-73-7-----	Fluorene	740	U
100-01-6-----	4-Nitroaniline	1800	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1800	U
86-30-6-----	N-Nitrosodiphenylamine (1)	740	U
101-55-3-----	4-Bromophenyl-phenylether	740	U
118-74-1-----	Hexachlorobenzene	740	U
87-86-5-----	Pentachlorophenol	1800	U
85-01-8-----	Phenanthrene	330	J
120-12-7-----	Anthracene	77	J
86-74-8-----	Carbazole	740	U
84-74-2-----	Di-n-Butylphthalate	740	U
206-44-0-----	Fluoranthene	790	
129-00-0-----	Pyrene	450	J
85-68-7-----	Butylbenzylphthalate	130	J
91-94-1-----	3,3'-Dichlorobenzidine	740	U
56-55-3-----	Benzo(a)Anthracene	420	J
218-01-9-----	Chrysene	340	J
117-81-7-----	bis(2-Ethylhexyl) Phthalate	350	J
117-84-0-----	Di-n-Octyl Phthalate	740	U
205-99-2-----	Benzo(b) Fluoranthene	610	JX
207-08-9-----	Benzo(k) Fluoranthene	610	JX
50-32-8-----	Benzo(a) Pyrene	360	J
193-39-5-----	Indeno(1,2,3-cd) Pyrene	220	J
53-70-3-----	Dibenz(a,h) Anthracene	78	J
191-24-2-----	Benzo(g,h,i) Perylene	190	J

(1) - Cannot be separated from Diphenylamine

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY14

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477484

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GR077484C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 56 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/11/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.2

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Number TICs found: 26

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.50	370	J
2.	TETRACHLOROETHANE	4.62	220	J
3.	UNKNOWN	4.72	1400	J
4.	ALDOL	4.85	1100	ABJ
5.	ALDOL	4.97	670	AJ
6. 100-52-7	BENZALDEHYDE	5.07	150	JN
7.	ALDOL	5.47	150	AJ
8.	UNKNOWN	5.95	2500	J
9.	UNKNOWN	7.10	300	J
10.	UNKNOWN	7.38	370	J
11.	BLANK CONTAMINANT	7.93	220	BJ
12.	BLANK CONTAMINANT	9.20	300	BJ
13.	UNKNOWN	12.32	220	J
14.	UNKNOWN CARBOXYLIC ACID	12.40	450	J
15.	UNKNOWN	13.42	220	J
16.	LABORATORY ARTIFACT	14.64	150	BJ
17.	UNKNOWN HYDROCARBON	15.15	1600	J
18.	UNKNOWN HYDROCARBON	16.09	2500	J
19.	UNKNOWN	16.54	670	J
20.	UNKNOWN	16.59	820	J
21.	UNKNOWN	16.84	1400	J
22.	UNKNOWN	17.24	13000	J
23.	BENZOFLUORANTHENE	17.37	600	J
24.	UNKNOWN	18.32	1700	J
25.	UNKNOWN	18.90	6800	J
26.	UNKNOWN	21.54	3400	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY14

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477484

Sample wt/vol: 30.10(g/ml)G

Lab File ID:

% Moisture: 56 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.2

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6-----	alpha-BHC	0.36	JP
319-85-7-----	beta-BHC	3.9	U
319-86-8-----	delta-BHC	3.9	U
58-89-9-----	gamma-BHC (Lindane)	40	E
76-44-8-----	Heptachlor	3.9	U
309-00-2-----	Aldrin	4.3	P
1024-57-3-----	Heptachlor epoxide	5.6	P
959-98-8-----	Endosulfan I	3.9	U
60-57-1-----	Dieldrin	7.6	U
72-55-9-----	4,4'-DDE	7.6	U
72-20-8-----	Endrin	4.4	JP
33213-65-9-----	Endosulfan II	7.6	U
72-54-8-----	4,4'-DDD	4.3	JP
1031-07-8-----	Endosulfan sulfate	0.69	JP
50-29-3-----	4,4'-DDT	1.5	JP
72-43-5-----	Methoxychlor	10.0	JPB
53494-70-5-----	Endrin ketone	4.6	JP
7421-93-4-----	Endrin aldehyde	3.1	JP
5103-71-9-----	alpha-Chlordane	6.2	P
5103-74-2-----	gamma-Chlordane	3.3	J
8001-35-2-----	Toxaphene	390	U
12674-11-2-----	Aroclor-1016	76	U
11104-28-2-----	Aroclor-1221	160	U
11141-16-5-----	Aroclor-1232	76	U
53469-21-9-----	Aroclor-1242	76	U
12672-29-6-----	Aroclor-1248	76	U
11097-69-1-----	Aroclor-1254	76	U
11096-82-5-----	Aroclor-1260	76	U

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY14DL

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477484 D20

Sample wt/vol: 30.10(g/ml)G

Lab File ID:

% Moisture: 56 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/13/92

Injection Volume: 2.0(uL)

Dilution Factor: 2

GPC Cleanup: (Y/N) Y

pH: 6.2

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
319-84-6	alpha-BHC	7.9	U
319-85-7	beta-BHC	7.9	U
319-86-8	delta-BHC	5.5	JPD
58-89-9	gamma-BHC (Lindane)	34	PD
76-44-8	Heptachlor	2.8	JPBD
309-00-2	Aldrin	3.2	JPD
1024-57-3	Heptachlor epoxide	1.9	JPD
959-98-8	Endosulfan I	14	PD
60-57-1	Dieldrin	3.3	JPD
72-55-9	4,4'-DDE	15	U
72-20-8	Endrin	4.5	JPD
33213-65-9	Endosulfan II	15	U
72-54-8	4,4'-DDD	3.8	JPD
1031-07-8	Endosulfan sulfate	15	U
50-29-3	4,4'-DDT	1.8	JPD
72-43-5	Methoxychlor	79	U
53494-70-5	Endrin ketone	15	U
7421-93-4	Endrin aldehyde	2.7	JPD
5103-71-9	alpha-Chlordane	7.9	U
5103-74-2	gamma-Chlordane	1.1	JPD
8001-35-2	Toxaphene	790	U
12674-11-2	Aroclor-1016	150	U
11104-28-2	Aroclor-1221	310	U
11141-16-5	Aroclor-1232	150	U
53469-21-9	Aroclor-1242	150	U
12672-29-6	Aroclor-1248	150	U
11097-69-1	Aroclor-1254	150	U
11096-82-5	Aroclor-1260	150	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY15

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477485

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077485B18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 19 Date Analyzed: 01/31/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:            (uL) Soil Aliquot Volume:            (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3-----	Chloromethane	12	U
74-83-9-----	Bromomethane	12	U
75-01-4-----	Vinyl Chloride	12	U
75-00-3-----	Chloroethane	12	U
75-09-2-----	Methylene Chloride	49	B
67-64-1-----	Acetone	61	B
75-15-0-----	Carbon Disulfide	12	U
75-35-4-----	1,1-Dichloroethene	12	U
75-34-3-----	1,1-Dichloroethane	12	U
540-59-0-----	1,2-Dichloroethene (total)	12	U
67-66-3-----	Chloroform	12	U
107-06-2-----	1,2-Dichloroethane	12	U
78-93-3-----	2-Butanone	12	U
71-55-6-----	1,1,1-Trichloroethane	12	U
56-23-5-----	Carbon Tetrachloride	12	U
75-27-4-----	Bromodichloromethane	12	U
78-87-5-----	1,2-Dichloropropane	12	U
10061-01-5-----	cis-1,3-Dichloropropene	12	U
79-01-6-----	Trichloroethene	12	U
124-48-1-----	Dibromochloromethane	12	U
79-00-5-----	1,1,2-Trichloroethane	12	U
71-43-2-----	Benzene	12	U
10061-02-6-----	Trans-1,3-Dichloropropene	12	U
75-25-2-----	Bromoform	12	U
108-10-1-----	4-Methyl-2-Pentanone	12	U
591-78-6-----	2-Hexanone	12	U
127-18-4-----	Tetrachloroethene	12	U
79-34-5-----	1,1,2,2-Tetrachloroethane	12	U
108-88-3-----	Toluene	12	U
108-90-7-----	Chlorobenzene	12	U
100-41-4-----	Ethylbenzene	12	U
100-42-5-----	Styrene	12	U
1330-20-7-----	Xylene (total)	12	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

096

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY15

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477485

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GRD77485A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 19 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 25.0

GPC Cleanup: (Y/N) Y pH: 6.7

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

108-95-2-----	Phenol	10000	U
111-44-4-----	bis(2-Chloroethyl) Ether	10000	U
95-57-8-----	2-Chlorophenol	10000	U
541-73-1-----	1,3-Dichlorobenzene	10000	U
106-46-7-----	1,4-Dichlorobenzene	10000	U
95-50-1-----	1,2-Dichlorobenzene	10000	U
95-48-7-----	2-Methylphenol	10000	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10000	U
106-44-5-----	4-Methylphenol	10000	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10000	U
67-72-1-----	Hexachloroethane	10000	U
98-95-3-----	Nitrobenzene	10000	U
78-59-1-----	Isophorone	10000	U
88-75-5-----	2-Nitrophenol	10000	U
105-67-9-----	2,4-Dimethylphenol	10000	U
111-91-1-----	bis(2-Chloroethoxy) Methane	10000	U
120-83-2-----	2,4-Dichlorophenol	10000	U
120-82-1-----	1,2,4-Trichlorobenzene	2300	J
91-20-3-----	Naphthalene	10000	U
106-47-8-----	4-Chloroaniline	10000	U
87-68-3-----	Hexachlorobutadiene	10000	U
59-50-7-----	4-Chloro-3-Methylphenol	10000	U
91-57-6-----	2-Methylnaphthalene	10000	U
77-47-4-----	Hexachlorocyclopentadiene	10000	U
88-06-2-----	2,4,6-Trichlorophenol	10000	U
95-95-4-----	2,4,5-Trichlorophenol	24000	U
91-58-7-----	2-Chloronaphthalene	10000	U
88-74-4-----	2-Nitroaniline	24000	U
131-11-3-----	Dimethyl Phthalate	10000	U
208-96-8-----	Acenaphthylene	10000	U
606-20-2-----	2,6-Dinitrotoluene	10000	U
99-09-2-----	3-Nitroaniline	24000	U
83-32-9-----	Acenaphthene	10000	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

714



1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY15

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477485

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GRD77485A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 19 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 25.0

PC Cleanup: (Y/N) Y pH: 6.7

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND Q

51-28-5-----	2,4-Dinitrophenol	24000	U
100-02-7-----	4-Nitrophenol	24000	U
132-64-9-----	Dibenzofuran	10000	U
121-14-2-----	2,4-Dinitrotoluene	10000	U
84-66-2-----	Diethylphthalate	10000	U
7005-72-3-----	4-Chlorophenyl-phenylether	10000	U
86-73-7-----	Fluorene	10000	U
100-01-6-----	4-Nitroaniline	24000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	24000	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10000	U
101-55-3-----	4-Bromophenyl-phenylether	10000	U
118-74-1-----	Hexachlorobenzene	10000	U
87-86-5-----	Pentachlorophenol	24000	U
85-01-8-----	Phenanthrene	2900	J
120-12-7-----	Anthracene	10000	U
86-74-8-----	Carbazole	10000	U
84-74-2-----	Di-n-Butylphthalate	10000	U
206-44-0-----	Fluoranthene	7100	J
129-00-0-----	Pyrene	4900	J
85-68-7-----	Butylbenzylphthalate	10000	U
91-94-1-----	3,3'-Dichlorobenzidine	10000	U
56-55-3-----	Benzo(a)Anthracene	3000	J
218-01-9-----	Chrysene	3800	J
117-81-7-----	bis(2-Ethylhexyl)Phthalate	1300	J
117-84-0-----	Di-n-Octyl Phthalate	10000	U
205-99-2-----	Benzo(b)Fluoranthene	8000	JX
207-08-9-----	Benzo(k)Fluoranthene	8000	JX
50-32-8-----	Benzo(a)Pyrene	3700	J
193-39-5-----	Indeno(1,2,3-cd)Pyrene	2400	J
53-70-3-----	Dibenz(a,h)Anthracene	10000	U
191-24-2-----	Benzo(g,h,i)Perylene	2100	J

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

715

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY15

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477485

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GRD77485A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 19 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 25.0

GPC Cleanup: (Y/N) Y pH: 6.7

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Number TICs found: 20

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	10.87	23000	J
2.	UNKNOWN	11.50	50000	J
3.	UNKNOWN	11.79	22000	J
4.	UNKNOWN	11.95	24000	J
5.	UNKNOWN	11.99	21000	J
6.	UNKNOWN	12.05	22000	J
7.	UNKNOWN	12.39	53000	J
8.	UNKNOWN	12.49	29000	J
9.	PCB	13.20	21000	J
10.	PCB	13.47	23000	J
11.	PCB	13.84	29000	J
12.	PCB	13.97	32000	J
13.	PCB	14.10	85000	J
14.	PCB	14.39	100000	J
15.	PCB	14.52	36000	J
16.	PCB	14.67	69000	J
17.	PCB	14.82	38000	J
18.	PCB	15.09	30000	J
19.	PCB	15.37	70000	J
20.	PCB	15.69	23000	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY15

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477485 R1D21

Sample wt/vol: 30.00(g/ml)G

Lab File ID: PR077485A22

% Moisture: 19 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/20/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/21/92

Injection Volume: 2.0(uL)

Dilution Factor: 20

GPC Cleanup: (Y/N) Y

pH: 6.7

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6	alpha-BHC	11	JP
319-85-7	beta-BHC	42	U
319-86-8	delta-BHC	42	U
58-89-9	gamma-BHC (Lindane)	3.4	JP
76-44-8	Heptachlor	42	U
309-00-2	Aldrin	12	JP
1024-57-3	Heptachlor epoxide	42	U
959-98-8	Endosulfan I	42	U
60-57-1	Dieldrin	81	U
72-55-9	4,4'-DDE	81	U
72-20-8	Endrin	81	U
33213-65-9	Endosulfan II	81	U
72-54-8	4,4'-DDD	81	U
1031-07-8	Endosulfan sulfate	81	U
50-29-3	4,4'-DDT	81	U
72-43-5	Methoxychlor	420	U
53494-70-5	Endrin ketone	81	U
7421-93-4	Endrin aldehyde	81	U
5103-71-9	alpha-Chlordane	42	U
5103-74-2	gamma-Chlordane	42	U
8001-35-2	Toxaphene	4200	U
12674-11-2	Aroclor-1016	810	U
11104-28-2	Aroclor-1221	1700	U
11141-16-5	Aroclor-1232	810	U
53469-21-9	Aroclor-1242	810	U
12672-29-6	Aroclor-1248	810	U
11097-69-1	Aroclor-1254	810	U
11096-82-5	Aroclor-1260	170000	PCB

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY15DL

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477485 R1D22

Sample wt/vol: 30.00(g/ml)G

Lab File ID: PR077485A22

% Moisture: 19 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/20/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/22/92

Injection Volume: 2.0(uL)

Dilution Factor: 200

GPC Cleanup: (Y/N) Y

pH: 6.7

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6-----	alpha-BHC	420	U
319-85-7-----	beta-BHC	420	U
319-86-8-----	delta-BHC	420	U
58-89-9-----	gamma-BHC (Lindane)	420	U
76-44-8-----	Heptachlor	420	U
309-00-2-----	Aldrin	420	U
1024-57-3-----	Heptachlor epoxide	420	U
959-98-8-----	Endosulfan I	420	U
60-57-1-----	Dieldrin	810	U
72-55-9-----	4,4'-DDE	810	U
72-20-8-----	Endrin	810	U
33213-65-9-----	Endosulfan II	810	U
72-54-8-----	4,4'-DDD	810	U
1031-07-8-----	Endosulfan sulfate	810	U
50-29-3-----	4,4'-DDT	810	U
72-43-5-----	Methoxychlor	4200	U
53494-70-5-----	Endrin ketone	810	U
7421-93-4-----	Endrin aldehyde	810	U
5103-71-9-----	alpha-Chlordane	420	U
5103-74-2-----	gamma-Chlordane	420	U
8001-35-2-----	Toxaphene	42000	U
12674-11-2-----	Aroclor-1016	8100	U
11104-28-2-----	Aroclor-1221	17000	U
11141-16-5-----	Aroclor-1232	8100	U
53469-21-9-----	Aroclor-1242	8100	U
12672-29-6-----	Aroclor-1248	8100	U
11097-69-1-----	Aroclor-1254	8100	U
11096-82-5-----	Aroclor-1260	300000	PCBD

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY16

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477486

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077486B18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 62 Date Analyzed: 01/31/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	26	U
74-83-9	-----Bromomethane	26	U
75-01-4	-----Vinyl Chloride	26	U
75-00-3	-----Chloroethane	26	U
75-09-2	-----Methylene Chloride	120	B
67-64-1	-----Acetone	420	B
75-15-0	-----Carbon Disulfide	26	U
75-35-4	-----1,1-Dichloroethene	26	U
75-34-3	-----1,1-Dichloroethane	26	U
540-59-0	-----1,2-Dichloroethene (total)	26	U
67-66-3	-----Chloroform	26	U
107-06-2	-----1,2-Dichloroethane	26	U
78-93-3	-----2-Butanone	160	
71-55-6	-----1,1,1-Trichloroethane	26	U
56-23-5	-----Carbon Tetrachloride	26	U
75-27-4	-----Bromodichloromethane	26	U
78-87-5	-----1,2-Dichloropropane	26	U
10061-01-5	-----cis-1,3-Dichloropropene	26	U
79-01-6	-----Trichloroethene	26	U
124-48-1	-----Dibromochloromethane	26	U
79-00-5	-----1,1,2-Trichloroethane	26	U
71-43-2	-----Benzene	26	U
10061-02-6	-----Trans-1,3-Dichloropropene	26	U
75-25-2	-----Bromoform	26	U
108-10-1	-----4-Methyl-2-Pentanone	340	
591-78-6	-----2-Hexanone	26	U
127-18-4	-----Tetrachloroethene	26	U
79-34-5	-----1,1,2,2-Tetrachloroethane	26	U
108-88-3	-----Toluene	120	
108-90-7	-----Chlorobenzene	26	U
100-41-4	-----Ethylbenzene	100	
100-42-5	-----Styrene	26	U
1330-20-7	-----Xylene (total)	690	

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY16

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477486

Sample wt/vol: 30.0 (g/mL) G Lab File ID: GRD77486B08

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 62 decanted: (Y/N) N Date Extracted: 02/16/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/18/92

Injection Volume: 2.0 (uL) Dilution Factor: 20.0

GPC Cleanup: (Y/N) Y pH: 6.0

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

108-95-2-----	Phenol	17000	U
111-44-4-----	bis(2-Chloroethyl) Ether	17000	U
95-57-8-----	2-Chlorophenol	17000	U
541-73-1-----	1,3-Dichlorobenzene	17000	U
106-46-7-----	1,4-Dichlorobenzene	17000	U
95-50-1-----	1,2-Dichlorobenzene	17000	U
95-48-7-----	2-Methylphenol	17000	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	17000	U
106-44-5-----	4-Methylphenol	17000	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	17000	U
67-72-1-----	Hexachloroethane	17000	U
98-95-3-----	Nitrobenzene	17000	U
78-59-1-----	Isophorone	17000	U
88-75-5-----	2-Nitrophenol	17000	U
105-67-9-----	2,4-Dimethylphenol	17000	U
111-91-1-----	bis(2-Chloroethoxy) Methane	17000	U
120-83-2-----	2,4-Dichlorophenol	17000	U
120-82-1-----	1,2,4-Trichlorobenzene	17000	U
91-20-3-----	Naphthalene	4500	J
106-47-8-----	4-Chloroaniline	17000	U
87-68-3-----	Hexachlorobutadiene	17000	U
59-50-7-----	4-Chloro-3-Methylphenol	17000	U
91-57-6-----	2-Methylnaphthalene	8000	J
77-47-4-----	Hexachlorocyclopentadiene	17000	U
88-06-2-----	2,4,6-Trichlorophenol	17000	U
95-95-4-----	2,4,5-Trichlorophenol	42000	U
91-58-7-----	2-Chloronaphthalene	17000	U
88-74-4-----	2-Nitroaniline	42000	U
131-11-3-----	Dimethyl Phthalate	17000	U
208-96-8-----	Acenaphthylene	17000	U
606-20-2-----	2,6-Dinitrotoluene	17000	U
99-09-2-----	3-Nitroaniline	42000	U
83-32-9-----	Acenaphthene	17000	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

773

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY16

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HO SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477486

Sample wt/vol: 30.0 (g/mL) G Lab File ID: GRD77486B08

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 62 decanted: (Y/N) N Date Extracted: 02/16/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/18/92

Injection Volume: 2.0 (uL) Dilution Factor: 20.0

GPC Cleanup: (Y/N) Y pH: 6.0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	Q
51-28-5-----	2,4-Dinitrophenol	42000 U
100-02-7-----	4-Nitrophenol	42000 U
132-64-9-----	Dibenzofuran	17000 U
121-14-2-----	2,4-Dinitrotoluene	17000 U
84-66-2-----	Diethylphthalate	17000 U
7005-72-3-----	4-Chlorophenyl-phenylether	17000 U
86-73-7-----	Fluorene	2600 J
100-01-6-----	4-Nitroaniline	42000 U
534-52-1-----	4,6-Dinitro-2-Methylphenol	42000 U
86-30-6-----	N-Nitrosodiphenylamine (1)	17000 U
101-55-3-----	4-Bromophenyl-phenylether	17000 U
118-74-1-----	Hexachlorobenzene	17000 U
87-86-5-----	Pentachlorophenol	42000 U
85-01-8-----	Phenanthrene	6500 J
120-12-7-----	Anthracene	17000 U
86-74-8-----	Carbazole	17000 U
84-74-2-----	Di-n-Butylphthalate	17000 U
206-44-0-----	Fluoranthene	4900 J
129-00-0-----	Pyrene	11000 J
85-68-7-----	Butylbenzylphthalate	17000 U
91-94-1-----	3,3'-Dichlorobenzidine	17000 U
56-55-3-----	Benzo(a)Anthracene	5200 J
218-01-9-----	Chrysene	4300 J
117-81-7-----	bis(2-Ethylhexyl) Phthalate	7500 J
117-84-0-----	Di-n-Octyl Phthalate	17000 U
205-99-2-----	Benzo(b) Fluoranthene	3900 J
207-08-9-----	Benzo(k) Fluoranthene	17000 U
50-32-8-----	Benzo(a) Pyrene	3700 J
193-39-5-----	Indeno(1,2,3-cd) Pyrene	17000 U
53-70-3-----	Dibenz(a,h) Anthracene	17000 U
191-24-2-----	Benzo(g,h,i) Perylene	7800 J

(1) - Cannot be separated from Diphenylamine

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY16

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477486

Sample wt/vol: 30.0 (g/mL) G Lab File ID: GRD77486B08

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 62 decanted: (Y/N) N Date Extracted: 02/16/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/18/92

Injection Volume: 2.0 (uL) Dilution Factor: 20.0

GPC Cleanup: (Y/N) Y pH: 6.0

Number TICs found: 20

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN HYDROCARBON	11.32	63000	J
2.	UNKNOWN HYDROCARBON	11.79	97000	J
3.	UNKNOWN HYDROCARBON	11.97	65000	J
4.	UNKNOWN HYDROCARBON	12.14	67000	J
5.	UNKNOWN HYDROCARBON	12.25	110000	J
6.	UNKNOWN HYDROCARBON	12.42	77000	J
7.	UNKNOWN HYDROCARBON	12.54	93000	J
8.	UNKNOWN	12.70	70000	J
9.	UNKNOWN HYDROCARBON	12.87	150000	J
10.	UNKNOWN	12.97	61000	J
11.	UNKNOWN HYDROCARBON	13.32	130000	J
12.	UNKNOWN CYCLIC HYDROCARBON	13.65	82000	J
13.	UNKNOWN HYDROCARBON	13.79	130000	J
14.	UNKNOWN HYDROCARBON	14.10	130000	J
15.	UNKNOWN HYDROCARBON	14.32	110000	J
16.	UNKNOWN HYDROCARBON	14.40	56000	J
17.	UNKNOWN HYDROCARBON	14.95	140000	J
18.	UNKNOWN	15.09	93000	J
19.	UNKNOWN	15.70	120000	J
20.	UNKNOWN	16.62	90000	J



1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY16

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477486

Sample wt/vol: 30.40(g/ml)G

Lab File ID:

% Moisture: 62 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) X

pH: 6.0

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
---------	----------	--	---

319-84-6-----	alpha-BHC	1.1	JP
319-85-7-----	beta-BHC	3.9	JP
319-86-8-----	delta-BHC	4.4	U
58-89-9-----	gamma-BHC (Lindane)	4.4	P
76-44-8-----	Heptachlor	0.16	JPB
309-00-2-----	Aldrin	2.6	JP
1024-57-3-----	Heptachlor epoxide	4.4	U
959-98-8-----	Endosulfan I	0.37	JP
60-57-1-----	Dieldrin	37	P
72-55-9-----	4,4'-DDE	13	P
72-20-8-----	Endrin	46	P
33213-65-9-----	Endosulfan II	64	P
72-54-8-----	4,4'-DDD	8.6	U
1031-07-8-----	Endosulfan sulfate	19	P
50-29-3-----	4,4'-DDT	8.6	U
72-43-5-----	Methoxychlor	180	PB
53494-70-5-----	Endrin ketone	8.6	U
7421-93-4-----	Endrin aldehyde	53	P
5103-71-9-----	alpha-Chlordane	11	P
5103-74-2-----	gamma-Chlordane	11	P
8001-35-2-----	Toxaphene	440	U
12674-11-2-----	Aroclor-1016	86	U
11104-28-2-----	Aroclor-1221	170	U
11141-16-5-----	Aroclor-1232	86	U
53469-21-9-----	Aroclor-1242	86	U
12672-29-6-----	Aroclor-1248	86	U
11097-69-1-----	Aroclor-1254	86	U
11096-82-5-----	Aroclor-1260	86	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY17

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477487

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077487A18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 22 Date Analyzed: 01/31/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	13	U
74-83-9-----	Bromomethane	13	U
75-01-4-----	Vinyl Chloride	13	U
75-00-3-----	Chloroethane	13	U
75-09-2-----	Methylene Chloride	45	B
67-64-1-----	Acetone	28	B
75-15-0-----	Carbon Disulfide	13	U
75-35-4-----	1,1-Dichloroethene	13	U
75-34-3-----	1,1-Dichloroethane	13	U
540-59-0-----	1,2-Dichloroethene (total)	13	U
67-66-3-----	Chloroform	13	U
107-06-2-----	1,2-Dichloroethane	13	U
78-93-3-----	2-Butanone	13	U
71-55-6-----	1,1,1-Trichloroethane	13	U
56-23-5-----	Carbon Tetrachloride	13	U
75-27-4-----	Bromodichloromethane	13	U
78-87-5-----	1,2-Dichloropropane	13	U
10061-01-5-----	cis-1,3-Dichloropropene	13	U
79-01-6-----	Trichloroethene	13	U
124-48-1-----	Dibromochloromethane	13	U
79-00-5-----	1,1,2-Trichloroethane	13	U
71-43-2-----	Benzene	13	U
10061-02-6-----	Trans-1,3-Dichloropropene	13	U
75-25-2-----	Bromoform	13	U
108-10-1-----	4-Methyl-2-Pentanone	13	U
591-78-6-----	2-Hexanone	13	U
127-18-4-----	Tetrachloroethene	13	U
79-34-5-----	1,1,2,2-Tetrachloroethane	13	U
108-88-3-----	Toluene	13	U
108-90-7-----	Chlorobenzene	13	U
100-41-4-----	Ethylbenzene	13	U
100-42-5-----	Styrene	13	U
1330-20-7-----	Xylene (total)	13	U

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY17

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477487

Sample wt/vol: 30.2 (g/mL) G Lab File ID: GRD77487A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 22 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/11/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

SPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	Q
108-95-2	Phenol	840 U
111-44-4	bis(2-Chloroethyl) Ether	840 U
95-57-8	2-Chlorophenol	840 U
541-73-1	1,3-Dichlorobenzene	840 U
106-46-7	1,4-Dichlorobenzene	840 U
95-50-1	1,2-Dichlorobenzene	840 U
95-48-7	2-Methylphenol	840 U
108-60-1	2,2'-Oxybis(1-Chloropropane)	840 U
106-44-5	4-Methylphenol	840 U
621-64-7	N-Nitroso-Di-n-Propylamine	840 U
67-72-1	Hexachloroethane	840 U
98-95-3	Nitrobenzene	840 U
78-59-1	Isophorone	840 U
88-75-5	2-Nitrophenol	840 U
105-67-9	2,4-Dimethylphenol	840 U
111-91-1	bis(2-Chloroethoxy) Methane	840 U
120-83-2	2,4-Dichlorophenol	840 U
120-82-1	1,2,4-Trichlorobenzene	840 U
91-20-3	Naphthalene	840 U
106-47-8	4-Chloroaniline	840 U
87-68-3	Hexachlorobutadiene	840 U
59-50-7	4-Chloro-3-Methylphenol	840 U
91-57-6	2-Methylnaphthalene	840 U
77-47-4	Hexachlorocyclopentadiene	840 U
88-06-2	2,4,6-Trichlorophenol	840 U
95-95-4	2,4,5-Trichlorophenol	2000 U
91-58-7	2-Chloronaphthalene	840 U
88-74-4	2-Nitroaniline	2000 U
131-11-3	Dimethyl Phthalate	840 U
208-96-8	Acenaphthylene	840 U
606-20-2	2,6-Dinitrotoluene	840 U
99-09-2	3-Nitroaniline	2000 U
83-32-9	Acenaphthene	98 J

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

833

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY17

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477487

Sample wt/vol: 30.2 (g/mL) G Lab File ID: GRD77487A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 22 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/11/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

GPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND Q

51-28-5-----	2,4-Dinitrophenol	2000	U
100-02-7-----	4-Nitrophenol	2000	U
132-64-9-----	Dibenzofuran	840	U
121-14-2-----	2,4-Dinitrotoluene	840	U
84-66-2-----	Diethylphthalate	840	U
7005-72-3-----	4-Chlorophenyl-phenylether	840	U
86-73-7-----	Fluorene	840	U
100-01-6-----	4-Nitroaniline	2000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	2000	U
86-30-6-----	N-Nitrosodiphenylamine (1)	840	U
101-55-3-----	4-Bromophenyl-phenylether	840	U
118-74-1-----	Hexachlorobenzene	840	U
87-86-5-----	Pentachlorophenol	2000	U
85-01-8-----	Phenanthrene	620	J
120-12-7-----	Anthracene	840	U
86-74-8-----	Carbazole	100	J
84-74-2-----	Di-n-Butylphthalate	840	U
206-44-0-----	Fluoranthene	1200	
129-00-0-----	Pyrene	920	
85-68-7-----	Butylbenzylphthalate	840	U
91-94-1-----	3,3'-Dichlorobenzidine	840	U
56-55-3-----	Benzo(a)Anthracene	630	J
218-01-9-----	Chrysene	610	J
117-81-7-----	bis(2-Ethylhexyl)Phthalate	840	U
117-84-0-----	Di-n-Octyl Phthalate	840	U
205-99-2-----	Benzo(b)Fluoranthene	1500	X
207-08-9-----	Benzo(k)Fluoranthene	1500	X
50-32-8-----	Benzo(a)Pyrene	480	J
193-39-5-----	Indeno(1,2,3-cd)Pyrene	600	J
53-70-3-----	Dibenz(a,h)Anthracene	200	J
191-24-2-----	Benzo(g,h,i)Perylene	520	J

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

834

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY17

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477487

Sample wt/vol: 30.2 (g/mL) G Lab File ID: GRD77487A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 22 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/11/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

PC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Number TICs found: 23

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.47	1500	J
2.	TETRACHLOROETHANE	4.58	420	J
3.	UNKNOWN	4.70	590	J
4.	ALDOL	4.82	1300	ABJ
5.	ALDOL	4.95	2900	AJ
6.	UNKNOWN	5.03	250	J
7.	UNKNOWN	5.27	680	J
8.	UNKNOWN	5.92	850	J
9.	THIENOPYRIDINE	7.57	1100	J
10. 615-22-5	BENZOTHAZOLE, 2-(METHYLTHIO	10.44	510	JN
11.	UNKNOWN	10.79	510	J
12.	UNKNOWN	12.44	340	J
13.	UNKNOWN	12.65	1400	J
14.	UNKNOWN	13.15	340	J
15.	UNKNOWN	13.90	420	J
16.	LABORATORY ARTIFACT	14.62	170	BJ
17.	UNKNOWN	17.20	1300	J
18.	UNKNOWN PAH	17.34	850	J
19.	UNKNOWN	18.80	340	J
20.	UNKNOWN	18.84	420	J
21.	UNKNOWN	18.89	170	J
22.	UNKNOWN	21.52	1300	J
23.	UNKNOWN	23.44	250	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY17

Lab Name: COMPUCHEM RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477487

Sample wt/vol: 30.40(g/ml)G

Lab File ID:

% Moisture: 22 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.6

Sulfur Cleanup: (Y/N) N

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

319-84-6-----	alpha-BHC	2.2	U
319-85-7-----	beta-BHC	2.2	U
319-86-8-----	delta-BHC	2.2	U
58-89-9-----	gamma-BHC (Lindane)	14	P
76-44-8-----	Heptachlor	16	PB
309-00-2-----	Aldrin	14	
1024-57-3-----	Heptachlor epoxide	0.19	JP
959-98-8-----	Endosulfan I	2.2	U
60-57-1-----	Dieldrin	31	P
72-55-9-----	4,4'-DDE	4.2	U
72-20-8-----	Endrin	39	P
33213-65-9-----	Endosulfan II	2.2	JP
72-54-8-----	4,4'-DDD	4.2	U
1031-07-8-----	Endosulfan sulfate	4.2	U
50-29-3-----	4,4'-DDT	31	P
72-43-5-----	Methoxychlor	31	PB
53494-70-5-----	Endrin ketone	4.2	U
7421-93-4-----	Endrin aldehyde	4.2	U
5103-71-9-----	alpha-Chlordane	2.2	U
5103-74-2-----	gamma-Chlordane	0.093	JP
8001-35-2-----	Toxaphene	220	U
12674-11-2-----	Aroclor-1016	42	U
11104-28-2-----	Aroclor-1221	85	U
11141-16-5-----	Aroclor-1232	42	U
53469-21-9-----	Aroclor-1242	42	U
12672-29-6-----	Aroclor-1248	42	U
11097-69-1-----	Aroclor-1254	42	U
11096-82-5-----	Aroclor-1260	42	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY18

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477489

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077489A18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 28 Date Analyzed: 01/31/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:            (uL) Soil Aliquot Volume:            (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	14	U
74-83-9-----	Bromomethane	14	U
75-01-4-----	Vinyl Chloride	14	U
75-00-3-----	Chloroethane	14	U
75-09-2-----	Methylene Chloride	52	B
67-64-1-----	Acetone	25	B
75-15-0-----	Carbon Disulfide	14	U
75-35-4-----	1,1-Dichloroethene	14	U
75-34-3-----	1,1-Dichloroethane	14	U
540-59-0-----	1,2-Dichloroethene (total)	14	U
67-66-3-----	Chloroform	14	U
107-06-2-----	1,2-Dichloroethane	14	U
78-93-3-----	2-Butanone	14	U
71-55-6-----	1,1,1-Trichloroethane	14	U
56-23-5-----	Carbon Tetrachloride	14	U
75-27-4-----	Bromodichloromethane	14	U
78-87-5-----	1,2-Dichloropropane	14	U
10061-01-5-----	cis-1,3-Dichloropropene	14	U
79-01-6-----	Trichloroethene	14	U
124-48-1-----	Dibromochloromethane	14	U
79-00-5-----	1,1,2-Trichloroethane	14	U
71-43-2-----	Benzene	14	U
10061-02-6-----	Trans-1,3-Dichloropropene	14	U
75-25-2-----	Bromoform	14	U
108-10-1-----	4-Methyl-2-Pentanone	14	U
591-78-6-----	2-Hexanone	14	U
127-18-4-----	Tetrachloroethene	14	U
79-34-5-----	1,1,2,2-Tetrachloroethane	14	U
108-88-3-----	Toluene	14	U
108-90-7-----	Chlorobenzene	14	U
100-41-4-----	Ethylbenzene	14	U
100-42-5-----	Styrene	14	U
1330-20-7-----	Xylene (total)	14	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

153

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY18

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477489

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GR077489C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 28 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/10/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.1

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND Q

108-95-2-----	Phenol	450	U
111-44-4-----	bis(2-Chloroethyl) Ether	450	U
95-57-8-----	2-Chlorophenol	450	U
541-73-1-----	1,3-Dichlorobenzene	450	U
106-46-7-----	1,4-Dichlorobenzene	450	U
95-50-1-----	1,2-Dichlorobenzene	450	U
95-48-7-----	2-Methylphenol	450	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	450	U
106-44-5-----	4-Methylphenol	450	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	450	U
67-72-1-----	Hexachloroethane	450	U
98-95-3-----	Nitrobenzene	450	U
78-59-1-----	Isophorone	450	U
88-75-5-----	2-Nitrophenol	450	U
105-67-9-----	2,4-Dimethylphenol	450	U
111-91-1-----	bis(2-Chloroethoxy) Methane	450	U
120-83-2-----	2,4-Dichlorophenol	450	U
120-82-1-----	1,2,4-Trichlorobenzene	450	U
91-20-3-----	Naphthalene	450	U
106-47-8-----	4-Chloroaniline	450	U
87-68-3-----	Hexachlorobutadiene	450	U
59-50-7-----	4-Chloro-3-Methylphenol	450	U
91-57-6-----	2-Methylnaphthalene	450	U
77-47-4-----	Hexachlorocyclopentadiene	450	U
88-06-2-----	2,4,6-Trichlorophenol	450	U
95-95-4-----	2,4,5-Trichlorophenol	1100	U
91-58-7-----	2-Chloronaphthalene	450	U
88-74-4-----	2-Nitroaniline	1100	U
131-11-3-----	Dimethyl Phthalate	450	U
208-96-8-----	Acenaphthylene	450	U
606-20-2-----	2,6-Dinitrotoluene	450	U
99-09-2-----	3-Nitroaniline	1100	U
83-32-9-----	Acenaphthene	89	J

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

895



1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY18

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477489

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GR077489C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 28 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/10/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.1

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS:	Q
51-28-5-----	2,4-Dinitrophenol	1100	U
100-02-7-----	4-Nitrophenol	1100	U
132-64-9-----	Dibenzofuran	450	U
121-14-2-----	2,4-Dinitrotoluene	450	U
84-66-2-----	Diethylphthalate	450	U
7005-72-3-----	4-Chlorophenyl-phenylether	450	U
86-73-7-----	Fluorene	66	J
100-01-6-----	4-Nitroaniline	1100	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1100	U
86-30-6-----	N-Nitrosodiphenylamine (1)	450	U
101-55-3-----	4-Bromophenyl-phenylether	450	U
118-74-1-----	Hexachlorobenzene	450	U
87-86-5-----	Pentachlorophenol	1100	U
85-01-8-----	Phenanthrene	1200	
120-12-7-----	Anthracene	140	J
86-74-8-----	Carbazole	180	J
84-74-2-----	Di-n-Butylphthalate	450	U
206-44-0-----	Fluoranthene	3700	E
129-00-0-----	Pyrene	2700	
85-68-7-----	Butylbenzylphthalate	450	U
91-94-1-----	3,3'-Dichlorobenzidine	450	U
56-55-3-----	Benzo(a) Anthracene	1700	
218-01-9-----	Chrysene	1600	
117-81-7-----	bis(2-Ethylhexyl) Phthalate	450	U
117-84-0-----	Di-n-Octyl Phthalate	450	U
205-99-2-----	Benzo(b) Fluoranthene	4600	EX
207-08-9-----	Benzo(k) Fluoranthene	4600	EX
50-32-8-----	Benzo(a) Pyrene	1200	
193-39-5-----	Indeno(1,2,3-cd) Pyrene	980	
53-70-3-----	Dibenz(a,h) Anthracene	350	J
191-24-2-----	Benzo(g,h,i) Perylene	950	

(1) - Cannot be separated from Diphenylamine

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY18

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477489

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GR077489C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 28 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/10/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.1

Number TICs found: 23

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.25	410	J
2.	UNKNOWN ALKENE	4.37	410	J
3.	UNKNOWN ALKENE	4.50	2200	J
4.	TETRACHLOROETHANE	4.62	320	J
5.	UNKNOWN	4.72	1500	J
6.	ALDOL	4.83	270	ABJ
7.	ALDOL	4.97	460	AJ
8.	UNKNOWN	5.07	370	J
9.	UNKNOWN	5.28	1100	J
10.	ALDOL	5.47	180	AJ
11.	UNKNOWN	5.93	1500	J
12.	UNKNOWN	7.40	180	J
13.	UNKNOWN	7.58	270	J
14.	UNKNOWN	12.32	180	J
15. 57-10-3	HEXADECANOIC ACID	12.40	230	JN
16.	UNKNOWN PAH	12.47	180	J
17.	UNKNOWN PNA	12.77	180	J
18.	UNKNOWN	16.07	460	J
19.	UNKNOWN HYDROCARBON	17.20	6900	J
20.	BENZOPYRENE	17.37	1500	J
21.	UNKNOWN	18.29	1400	J
22.	UNKNOWN	18.85	9100	J
23.	UNKNOWN	21.47	4000	J

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY18DL

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477489

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GDJ77489A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 28 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

IPC Cleanup: (Y/N) Y pH: 6.1

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

108-95-2-----	Phenol	900	U
111-44-4-----	bis(2-Chloroethyl) Ether	900	U
95-57-8-----	2-Chlorophenol	900	U
541-73-1-----	1,3-Dichlorobenzene	900	U
106-46-7-----	1,4-Dichlorobenzene	900	U
95-50-1-----	1,2-Dichlorobenzene	900	U
95-48-7-----	2-Methylphenol	900	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	900	U
106-44-5-----	4-Methylphenol	900	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	900	U
67-72-1-----	Hexachloroethane	900	U
98-95-3-----	Nitrobenzene	900	U
78-59-1-----	Isophorone	900	U
88-75-5-----	2-Nitrophenol	900	U
105-67-9-----	2,4-Dimethylphenol	900	U
111-91-1-----	bis(2-Chloroethoxy)Methane	900	U
120-83-2-----	2,4-Dichlorophenol	900	U
120-82-1-----	1,2,4-Trichlorobenzene	900	U
91-20-3-----	Naphthalene	900	U
106-47-8-----	4-Chloroaniline	900	U
87-68-3-----	Hexachlorobutadiene	900	U
59-50-7-----	4-Chloro-3-Methylphenol	900	U
91-57-6-----	2-Methylnaphthalene	900	U
77-47-4-----	Hexachlorocyclopentadiene	900	U
88-06-2-----	2,4,6-Trichlorophenol	900	U
95-95-4-----	2,4,5-Trichlorophenol	2200	U
91-58-7-----	2-Chloronaphthalene	900	U
88-74-4-----	2-Nitroaniline	2200	U
131-11-3-----	Dimethyl Phthalate	900	U
208-96-8-----	Acenaphthylene	900	U
606-20-2-----	2,6-Dinitrotoluene	900	U
99-09-2-----	3-Nitroaniline	2200	U
83-32-9-----	Acenaphthene	110	DJ

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

960

ORIGINAL

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY18DL

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477489

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GDJ77489A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 28 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

GPC Cleanup: (Y/N) Y pH: 6.1

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND		Q
51-28-5-----	2,4-Dinitrophenol	2200	U
100-02-7-----	4-Nitrophenol	2200	U
132-64-9-----	Dibenzofuran	900	U
121-14-2-----	2,4-Dinitrotoluene	900	U
84-66-2-----	Diethylphthalate	390	DJ
7005-72-3-----	4-Chlorophenyl-phenylether	900	U
86-73-7-----	Fluorene	92	DJ
100-01-6-----	4-Nitroaniline	2200	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	2200	U
86-30-6-----	N-Nitrosodiphenylamine (1)	900	U
101-55-3-----	4-Bromophenyl-phenylether	900	U
118-74-1-----	Hexachlorobenzene	900	U
87-86-5-----	Pentachlorophenol	2200	U
85-01-8-----	Phenanthrene	1400	D
120-12-7-----	Anthracene	240	DJ
86-74-8-----	Carbazole	260	DJ
84-74-2-----	Di-n-Butylphthalate	900	U
206-44-0-----	Fluoranthene	3900	D
129-00-0-----	Pyrene	3300	D
85-68-7-----	Butylbenzylphthalate	900	U
91-94-1-----	3,3'-Dichlorobenzidine	900	U
56-55-3-----	Benzo(a)Anthracene	1700	D
218-01-9-----	Chrysene	2100	D
117-81-7-----	bis(2-Ethylhexyl)Phthalate	900	U
117-84-0-----	Di-n-Octyl Phthalate	900	U
205-99-2-----	Benzo(b)Fluoranthene	4400	DX
207-08-9-----	Benzo(k)Fluoranthene	4400	DX
50-32-8-----	Benzo(a)Pyrene	1900	D
193-39-5-----	Indeno(1,2,3-cd)Pyrene	1700	D
53-70-3-----	Dibenz(a,h)Anthracene	590	DJ
191-24-2-----	Benzo(g,h,i)Perylene	1800	D

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

961

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY18DL

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477489

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GDJ77489A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 28 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

IPC Cleanup: (Y/N) Y pH: 6.1

Number TICs found: 26

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN ALKENE	4.37	1200	J
2.	TETRACHLOROETHANE	4.48	370	J
3.	UNKNOWN	4.58	1900	J
4.	ALDOL	4.72	820	ABJ
5.	UNKNOWN	4.95	370	J
6.	UNKNOWN	5.17	1300	J
7.	UNKNOWN	5.82	1600	J
8.	UNKNOWN	7.47	370	J
9.	BLANK CONTAMINANT	7.80	820	BJ
10.	BLANK CONTAMINANT	9.09	910	BJ
11.	UNKNOWN SUBST. PROPANOIC ACI	10.05	1700	J
12.	BLANK CONTAMINANT	11.17	2300	J
13.	UNKNOWN SILOXANE	12.02	2300	BJ
14. 57-10-3	HEXADECANOIC ACID	12.29	370	JN
15.	UNKNOWN	12.55	640	J
16.	UNKNOWN PNA	12.64	270	J
17.	UNKNOWN SILOXANE	12.80	2000	J
18.	UNKNOWN SILOXANE	14.17	1800	J
19.	LAB. ARTIFACT	14.50	370	BJ
20.	UNKNOWN SILOXANE	14.79	1600	J
21.	UNKNOWN SILOXANE	15.40	1300	J
22.	UNKNOWN HYDROCARBON	15.95	550	J
23.	UNKNOWN SILOXANE	16.02	1000	J
24.	UNKNOWN HYDROCARBON	17.05	5400	J
25.	UNKNOWN PAH	17.17	2700	J
26.	UNKNOWN	18.67	7200	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY18

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477489 R1

Sample wt/vol: 30.00(g/ml)G

Lab File ID:

% Moisture: 28 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/16/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/18/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.1

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG Q
319-84-6	alpha-BHC	2.4	U
319-85-7	beta-BHC	2.4	U
319-86-8	delta-BHC	2.4	U
58-89-9	gamma-BHC (Lindane)	2.4	U
76-44-8	Heptachlor	2.4	U
309-00-2	Aldrin	0.36	JP
1024-57-3	Heptachlor epoxide	0.54	JP
959-98-8	Endosulfan I	2.4	U
60-57-1	Dieldrin	0.27	JP
72-55-9	4,4'-DDE	2.8	J
72-20-8	Endrin	2.6	JPB
33213-65-9	Endosulfan II	2.6	J
72-54-8	4,4'-DDD	4.6	U
1031-07-8	Endosulfan sulfate	4.6	U
50-29-3	4,4'-DDT	0.48	JP
72-43-5	Methoxychlor	18	JPB
53494-70-5	Endrin ketone	4.6	U
7421-93-4	Endrin aldehyde	4.6	U
5103-71-9	alpha-Chlordane	2.4	U
5103-74-2	gamma-Chlordane	0.23	JP
8001-35-2	Toxaphene	240	U
12674-11-2	Aroclor-1016	46	U
11104-28-2	Aroclor-1221	93	U
11141-16-5	Aroclor-1232	46	U
53469-21-9	Aroclor-1242	46	U
12672-29-6	Aroclor-1248	46	U
11097-69-1	Aroclor-1254	46	U
11096-82-5	Aroclor-1260	46	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY19

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477490

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077490A18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 24 Date Analyzed: 01/31/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	13	U
74-83-9-----	Bromomethane	13	U
75-01-4-----	Vinyl Chloride	13	U
75-00-3-----	Chloroethane	13	U
75-09-2-----	Methylene Chloride	47	B
67-64-1-----	Acetone	57	B
75-15-0-----	Carbon Disulfide	13	U
75-35-4-----	1,1-Dichloroethene	13	U
75-34-3-----	1,1-Dichloroethane	13	U
540-59-0-----	1,2-Dichloroethene (total)	13	U
67-66-3-----	Chloroform	13	U
107-06-2-----	1,2-Dichloroethane	13	U
78-93-3-----	2-Butanone	13	U
71-55-6-----	1,1,1-Trichloroethane	13	U
56-23-5-----	Carbon Tetrachloride	13	U
75-27-4-----	Bromodichloromethane	13	U
78-87-5-----	1,2-Dichloropropane	13	U
10061-01-5-----	cis-1,3-Dichloropropene	13	U
79-01-6-----	Trichloroethene	13	U
124-48-1-----	Dibromochloromethane	13	U
79-00-5-----	1,1,2-Trichloroethane	13	U
71-43-2-----	Benzene	13	U
10061-02-6-----	Trans-1,3-Dichloropropene	13	U
75-25-2-----	Bromoform	13	U
108-10-1-----	4-Methyl-2-Pentanone	13	U
591-78-6-----	2-Hexanone	13	U
127-18-4-----	Tetrachloroethene	13	U
79-34-5-----	1,1,2,2-Tetrachloroethane	13	U
108-88-3-----	Toluene	13	U
108-90-7-----	Chlorobenzene	13	U
100-41-4-----	Ethylbenzene	13	U
100-42-5-----	Styrene	13	U
1330-20-7-----	Xylene (total)	13	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

165

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY19

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477490

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GRD77490A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 24 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

GPC Cleanup: (Y/N) Y pH: 6.2

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	Q
108-95-2-----	Phenol	860 U
111-44-4-----	bis(2-Chloroethyl) Ether	860 U
95-57-8-----	2-Chlorophenol	860 U
541-73-1-----	1,3-Dichlorobenzene	860 U
106-46-7-----	1,4-Dichlorobenzene	860 U
95-50-1-----	1,2-Dichlorobenzene	860 U
95-48-7-----	2-Methylphenol	860 U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	860 U
106-44-5-----	4-Methylphenol	860 U
621-64-7-----	N-Nitroso-Di-n-Propylamine	860 U
67-72-1-----	Hexachloroethane	860 U
98-95-3-----	Nitrobenzene	860 U
78-59-1-----	Isophorone	860 U
88-75-5-----	2-Nitrophenol	860 U
105-67-9-----	2,4-Dimethylphenol	860 U
111-91-1-----	bis(2-Chloroethoxy) Methane	860 U
120-83-2-----	2,4-Dichlorophenol	860 U
120-82-1-----	1,2,4-Trichlorobenzene	860 U
91-20-3-----	Naphthalene	150 J
106-47-8-----	4-Chloroaniline	860 U
87-68-3-----	Hexachlorobutadiene	860 U
59-50-7-----	4-Chloro-3-Methylphenol	860 U
91-57-6-----	2-Methylnaphthalene	95 J
77-47-4-----	Hexachlorocyclopentadiene	860 U
88-06-2-----	2,4,6-Trichlorophenol	860 U
95-95-4-----	2,4,5-Trichlorophenol	2100 U
91-58-7-----	2-Chloronaphthalene	860 U
88-74-4-----	2-Nitroaniline	2100 U
131-11-3-----	Dimethyl Phthalate	860 U
208-96-8-----	Acenaphthylene	860 U
606-20-2-----	2,6-Dinitrotoluene	860 U
99-09-2-----	3-Nitroaniline	2100 U
83-32-9-----	Acenaphthene	440 J

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1036



1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY19

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477490

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GRD77490A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 24 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 2.0

GPC Cleanup: (Y/N) Y pH: 6.2

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

51-28-5-----	2,4-Dinitrophenol	2100	U
100-02-7-----	4-Nitrophenol	2100	U
132-64-9-----	Dibenzofuran	130	J
121-14-2-----	2,4-Dinitrotoluene	860	U
84-66-2-----	Diethylphthalate	860	U
7005-72-3-----	4-Chlorophenyl-phenylether	860	U
86-73-7-----	Fluorene	180	J
100-01-6-----	4-Nitroaniline	2100	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	2100	U
86-30-6-----	N-Nitrosodiphenylamine (1)	860	U
101-55-3-----	4-Bromophenyl-phenylether	860	U
118-74-1-----	Hexachlorobenzene	860	U
87-86-5-----	Pentachlorophenol	2100	U
85-01-8-----	Phenanthrene	2300	
120-12-7-----	Anthracene	450	J
86-74-8-----	Carbazole	330	J
84-74-2-----	Di-n-Butylphthalate	860	U
206-44-0-----	Fluoranthene	4100	
129-00-0-----	Pyrene	2000	
85-68-7-----	Butylbenzylphthalate	860	U
91-94-1-----	3,3'-Dichlorobenzidine	860	U
56-55-3-----	Benzo(a)Anthracene	1400	
218-01-9-----	Chrysene	1500	
117-81-7-----	bis(2-Ethylhexyl) Phthalate	860	U
117-84-0-----	Di-n-Octyl Phthalate	860	U
205-99-2-----	Benzo(b) Fluoranthene	3000	X
207-08-9-----	Benzo(k) Fluoranthene	3000	X
50-32-8-----	Benzo(a) Pyrene	1600	
193-39-5-----	Indeno(1,2,3-cd) Pyrene	1100	
53-70-3-----	Dibenz(a,h) Anthracene	400	J
191-24-2-----	Benzo(g,h,i) Perylene	1000	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1037

ORIGINAL  
Red

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY19

Lab Name: COMPUCHEM RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
Matrix: (soil/water) SOIL Lab Sample ID: 477490  
Sample wt/vol: 30.4 (g/mL) G Lab File ID: GRD77490A21  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: 24 decanted: (Y/N) N Date Extracted: 02/04/92  
Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92  
Injection Volume: 2.0 (uL) Dilution Factor: 2.0  
GPC Cleanup: (Y/N) Y pH: 6.2

Number TICs found: 22 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.18	430	J
2.	UNKNOWN ALKENE	4.30	690	J
3.	UNKNOWN	4.45	5500	J
4.	TETRACHLOROETHANE	4.55	610	J
5.	UNKNOWN	4.65	2900	J
6.	ALDOL	4.80	1000	ABJ
7.	UNKNOWN	4.95	350	J
8.	UNKNOWN	5.02	610	J
9.	UNKNOWN	5.25	2400	J
10.	UNKNOWN	5.90	1200	J
11.	BLANK CONTAMINANT	11.27	260	BJ
12.	UNKNOWN PAH	12.45	350	J
13.	UNKNOWN PNA	12.75	350	J
14.	UNKNOWN	17.19	1000	J
15.	UNKNOWN PAH	17.32	1300	J
16.	UNKNOWN	18.27	350	J
17.	UNKNOWN	18.72	350	J
18.	UNKNOWN	18.77	430	J
19.	UNKNOWN	18.84	350	J
20.	UNKNOWN	21.47	780	J
21.	UNKNOWN	21.54	870	J
22.	UNKNOWN	23.45	610	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY19

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477490

Sample wt/vol: 30.30(g/ml)G

Lab File ID:

% Moisture: 24 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.2

Sulfur Cleanup: (Y/N) N

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

319-84-6-----	alpha-BHC	0.081	JP
319-85-7-----	beta-BHC	2.2	U
319-86-8-----	delta-BHC	2.2	U
58-89-9-----	gamma-BHC (Lindane)	2.2	U
76-44-8-----	Heptachlor	0.15	JPB
309-00-2-----	Aldrin	0.39	JP
1024-57-3-----	Heptachlor epoxide	0.41	JP
959-98-8-----	Endosulfan I	2.2	U
60-57-1-----	Dieldrin	0.10	JP
72-55-9-----	4,4'-DDE	1.1	J
72-20-8-----	Endrin	1.5	JP
33213-65-9-----	Endosulfan II	0.48	JP
72-54-8-----	4,4'-DDD	4.3	U
1031-07-8-----	Endosulfan sulfate	4.3	U
50-29-3-----	4,4'-DDT	4.3	U
72-43-5-----	Methoxychlor	6.5	JPB
53494-70-5-----	Endrin ketone	4.3	U
7421-93-4-----	Endrin aldehyde	4.3	U
5103-71-9-----	alpha-Chlordane	2.2	U
5103-74-2-----	gamma-Chlordane	2.2	U
8001-35-2-----	Toxaphene	220	U
12674-11-2-----	Aroclor-1016	43	U
11104-28-2-----	Aroclor-1221	87	U
11141-16-5-----	Aroclor-1232	43	U
53469-21-9-----	Aroclor-1242	43	U
12672-29-6-----	Aroclor-1248	43	U
11097-69-1-----	Aroclor-1254	43	U
11096-82-5-----	Aroclor-1260	43	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY20

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477491

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077491A18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 35 Date Analyzed: 01/31/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:            (uL) Soil Aliquot Volume:            (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	15	U
74-83-9-----	Bromomethane	15	U
75-01-4-----	Vinyl Chloride	15	U
75-00-3-----	Chloroethane	15	U
75-09-2-----	Methylene Chloride	70	B
67-64-1-----	Acetone	69	B
75-15-0-----	Carbon Disulfide	15	U
75-35-4-----	1,1-Dichloroethene	15	U
75-34-3-----	1,1-Dichloroethane	15	U
540-59-0-----	1,2-Dichloroethene (total)	15	U
67-66-3-----	Chloroform	15	U
107-06-2-----	1,2-Dichloroethane	15	U
78-93-3-----	2-Butanone	15	U
71-55-6-----	1,1,1-Trichloroethane	15	U
56-23-5-----	Carbon Tetrachloride	15	U
75-27-4-----	Bromodichloromethane	15	U
78-87-5-----	1,2-Dichloropropane	15	U
10061-01-5-----	cis-1,3-Dichloropropene	15	U
79-01-6-----	Trichloroethene	15	U
124-48-1-----	Dibromochloromethane	15	U
79-00-5-----	1,1,2-Trichloroethane	15	U
71-43-2-----	Benzene	15	U
10061-02-6-----	Trans-1,3-Dichloropropene	15	U
75-25-2-----	Bromoform	15	U
108-10-1-----	4-Methyl-2-Pentanone	15	U
591-78-6-----	2-Hexanone	15	U
127-18-4-----	Tetrachloroethene	15	U
79-34-5-----	1,1,2,2-Tetrachloroethane	15	U
108-88-3-----	Toluene	15	U
108-90-7-----	Chlorobenzene	15	U
100-41-4-----	Ethylbenzene	15	U
100-42-5-----	Styrene	15	U
1330-20-7-----	Xylene (total)	15	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

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1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY20

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477491

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GRD77491C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 35 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/11/92

Injection Volume: 2.0 (uL) Dilution Factor: 15.0

GPC Cleanup: (Y/N) Y pH: 6.7

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

108-95-2-----	Phenol	7500	U
111-44-4-----	bis(2-Chloroethyl) Ether	7500	U
95-57-8-----	2-Chlorophenol	7500	U
541-73-1-----	1,3-Dichlorobenzene	7500	U
106-46-7-----	1,4-Dichlorobenzene	7500	U
95-50-1-----	1,2-Dichlorobenzene	7500	U
95-48-7-----	2-Methylphenol	7500	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	7500	U
106-44-5-----	4-Methylphenol	7500	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	7500	U
67-72-1-----	Hexachloroethane	7500	U
98-95-3-----	Nitrobenzene	7500	U
78-59-1-----	Isophorone	7500	U
88-75-5-----	2-Nitrophenol	7500	U
105-67-9-----	2,4-Dimethylphenol	7500	U
111-91-1-----	bis(2-Chloroethoxy) Methane	7500	U
120-83-2-----	2,4-Dichlorophenol	7500	U
120-82-1-----	1,2,4-Trichlorobenzene	7500	U
91-20-3-----	Naphthalene	2500	J
106-47-8-----	4-Chloroaniline	7500	U
87-68-3-----	Hexachlorobutadiene	7500	U
59-50-7-----	4-Chloro-3-Methylphenol	7500	U
91-57-6-----	2-Methylnaphthalene	790	J
77-47-4-----	Hexachlorocyclopentadiene	7500	U
88-06-2-----	2,4,6-Trichlorophenol	7500	U
95-95-4-----	2,4,5-Trichlorophenol	18000	U
91-58-7-----	2-Chloronaphthalene	7500	U
88-74-4-----	2-Nitroaniline	18000	U
131-11-3-----	Dimethyl Phthalate	7500	U
208-96-8-----	Acenaphthylene	7500	U
606-20-2-----	2,6-Dinitrotoluene	7500	U
99-09-2-----	3-Nitroaniline	18000	U
83-32-9-----	Acenaphthene	9000	

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1108

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY20

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
 Matrix: (soil/water) SOIL Lab Sample ID: 477491  
 Sample wt/vol: 30.5 (g/mL) G Lab File ID: GRD77491C21  
 Level: (low/med) LOW Date Received: 01/29/92  
 % Moisture: 35 decanted: (Y/N) N Date Extracted: 02/04/92  
 Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/11/92  
 Injection Volume: 2.0 (uL) Dilution Factor: 15.0  
 GPC Cleanup: (Y/N) Y pH: 6.7

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
51-28-5-----	2,4-Dinitrophenol	18000	U
100-02-7-----	4-Nitrophenol	18000	U
132-64-9-----	Dibenzofuran	2900	J
121-14-2-----	2,4-Dinitrotoluene	7500	U
84-66-2-----	Diethylphthalate	7500	U
7005-72-3-----	4-Chlorophenyl-phenylether	7500	U
86-73-7-----	Fluorene	4100	J
100-01-6-----	4-Nitroaniline	18000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	18000	U
86-30-6-----	N-Nitrosodiphenylamine (1)	7500	U
101-55-3-----	4-Bromophenyl-phenylether	7500	U
118-74-1-----	Hexachlorobenzene	7500	U
87-86-5-----	Pentachlorophenol	18000	U
85-01-8-----	Phenanthrene	75000	E
120-12-7-----	Anthracene	14000	
86-74-8-----	Carbazole	11000	
84-74-2-----	Di-n-Butylphthalate	7500	U
206-44-0-----	Fluoranthene	110000	E
129-00-0-----	Pyrene	51000	
85-68-7-----	Butylbenzylphthalate	7500	U
91-94-1-----	3,3'-Dichlorobenzidine	7500	U
56-55-3-----	Benzo(a)Anthracene	33000	
218-01-9-----	Chrysene	32000	
117-81-7-----	bis(2-Ethylhexyl)Phthalate	1600	J
117-84-0-----	Di-n-Octyl Phthalate	7500	U
205-99-2-----	Benzo(b)Fluoranthene	65000	EX
207-08-9-----	Benzo(k)Fluoranthene	65000	EX
50-32-8-----	Benzo(a)Pyrene	30000	
193-39-5-----	Indeno(1,2,3-cd)Pyrene	11000	
53-70-3-----	Dibenz(a,h)Anthracene	5300	J
191-24-2-----	Benzo(g,h,i)Perylene	7600	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1109

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY20

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477491

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GRD77491C21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 35 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/11/92

Injection Volume: 2.0 (uL) Dilution Factor: 15.0

GPC Cleanup: (Y/N) Y pH: 6.7

Number TICs found: 22

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	ALDOL	4.85	3800	ABJ
2.	BLANK CONTAMINANT	9.22	2300	BJ
3.	HYDROXYNAPHTHALENEDIONE	9.79	23000	J
4.	UNKNOWN PAH	12.50	7600	J
5.	UNKNOWN PAH	13.82	7600	J
6.	UNKNOWN PAH	13.95	8300	J
7.	BENZO FLUORENE	14.12	18000	J
8.	BENZOFLUORENE	14.20	11000	J
9.	METHYL PYRENE	14.25	8300	J
10.	UNKNOWN PAH	14.29	7600	J
11.	UNKNOWN PAH	14.82	8300	J
12.	UNKNOWN PAH	14.99	12000	J
13.	UNKNOWN PAH	15.02	7600	J
14.	UNKNOWN PAH	15.05	13000	J
15.	UNKNOWN PAH	15.12	7600	J
16.	UNKNOWN PAH	15.49	12000	J
17.	UNKNOWN PAH	15.64	9800	J
18.	UNKNOWN PAH	15.69	7600	J
19.	UNKNOWN PAH	15.89	15000	J
20.	UNKNOWN	16.10	17000	J
21.	BENZOFLUORANTHENE	17.09	6100	J
22.	BENZOFLUORANTHENE	17.40	22000	J

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY20DL

Lab Name: COMPUCHEM.RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
 Matrix: (soil/water) SOIL Lab Sample ID: 477491  
 Sample wt/vol: 30.5 (g/mL) G Lab File ID: GDJ77491A21  
 Level: (low/med) LOW Date Received: 01/29/92  
 % Moisture: 35 decanted: (Y/N) N Date Extracted: 02/04/92  
 Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/13/92  
 Injection Volume: 2.0 (uL) Dilution Factor: 30.0  
 GPC Cleanup: (Y/N) Y pH: 6.7

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.

COMPOUND

Q

108-95-2-----	Phenol	15000	U
111-44-4-----	bis(2-Chloroethyl) Ether	15000	U
95-57-8-----	2-Chlorophenol	15000	U
541-73-1-----	1,3-Dichlorobenzene	15000	U
106-46-7-----	1,4-Dichlorobenzene	15000	U
95-50-1-----	1,2-Dichlorobenzene	15000	U
95-48-7-----	2-Methylphenol	15000	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	15000	U
106-44-5-----	4-Methylphenol	15000	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	15000	U
67-72-1-----	Hexachloroethane	15000	U
98-95-3-----	Nitrobenzene	15000	U
78-59-1-----	Isophorone	15000	U
88-75-5-----	2-Nitrophenol	15000	U
105-67-9-----	2,4-Dimethylphenol	15000	U
111-91-1-----	bis(2-Chloroethoxy)Methane	15000	U
120-83-2-----	2,4-Dichlorophenol	15000	U
120-82-1-----	1,2,4-Trichlorobenzene	15000	U
91-20-3-----	Naphthalene	2700	DJ
106-47-8-----	4-Chloroaniline	15000	U
87-68-3-----	Hexachlorobutadiene	15000	U
59-50-7-----	4-Chloro-3-Methylphenol	15000	U
91-57-6-----	2-Methylnaphthalene	15000	U
77-47-4-----	Hexachlorocyclopentadiene	15000	U
88-06-2-----	2,4,6-Trichlorophenol	15000	U
95-95-4-----	2,4,5-Trichlorophenol	36000	U
91-58-7-----	2-Chloronaphthalene	15000	U
88-74-4-----	2-Nitroaniline	36000	U
131-11-3-----	Dimethyl Phthalate	15000	U
208-96-8-----	Acenaphthylene	15000	U
606-20-2-----	2,6-Dinitrotoluene	15000	U
99-09-2-----	3-Nitroaniline	36000	U
83-32-9-----	Acenaphthene	10000	DJ

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1181



1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY20DL

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477491

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GDJ77491A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 35 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/13/92

Injection Volume: 2.0 (uL) Dilution Factor: 30.0

PC Cleanup: (Y/N) Y pH: 6.7

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	Q
51-28-5-----	2,4-Dinitrophenol	U
100-02-7-----	4-Nitrophenol	U
132-64-9-----	Dibenzofuran	DJ
121-14-2-----	2,4-Dinitrotoluene	U
84-66-2-----	Diethylphthalate	DJ
7005-72-3-----	4-Chlorophenyl-phenylether	U
86-73-7-----	Fluorene	DJ
100-01-6-----	4-Nitroaniline	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	U
86-30-6-----	N-Nitrosodiphenylamine (1)	U
101-55-3-----	4-Bromophenyl-phenylether	U
118-74-1-----	Hexachlorobenzene	U
87-86-5-----	Pentachlorophenol	U
85-01-8-----	Phenanthrene	D
120-12-7-----	Anthracene	DJ
86-74-8-----	Carbazole	DJ
84-74-2-----	Di-n-Butylphthalate	U
206-44-0-----	Fluoranthene	D
129-00-0-----	Pyrene	D
85-68-7-----	Butylbenzylphthalate	U
91-94-1-----	3,3'-Dichlorobenzidine	U
56-55-3-----	Benzo(a) Anthracene	D
218-01-9-----	Chrysene	D
117-81-7-----	bis(2-Ethylhexyl) Phthalate	DJ
117-84-0-----	Di-n-Octyl Phthalate	U
205-99-2-----	Benzo(b) Fluoranthene	DX
207-08-9-----	Benzo(k) Fluoranthene	DX
50-32-8-----	Benzo(a) Pyrene	D
193-39-5-----	Indeno(1,2,3-cd) Pyrene	D
53-70-3-----	Dibenz(a,h) Anthracene	DJ
191-24-2-----	Benzo(g,h,i) Perylene	DJ

(1) - Cannot be separated from Diphenylamine

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY20DL

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477491

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GDJ77491A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 35 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/13/92

Injection Volume: 2.0 (uL) Dilution Factor: 30.0

GPC Cleanup: (Y/N) Y pH: 6.7

Number TICs found: 24

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	BLANK CONTAMINANT	7.87	6100	BJ
2.	BLANK CONTAMINANT	9.14	7600	BJ
3.	HYDROXY NAPHTHALENEDIONE	9.69	20000	J
4.	BLANK CONTAMINANT	11.20	24000	BJ
5.	UNKNOWN SILOXANE	12.05	27000	J
6.	UNKNOWN PAH	12.27	4500	J
7.	UNKNOWN PAH	12.37	12000	J
8.	UNKNOWN	12.59	7600	J
9.	UNKNOWN SILOXANE	12.84	27000	J
10.	UNKNOWN PAH	13.99	6100	J
11.	UNKNOWN PAH	14.09	4500	J
12.	UNKNOWN PAH	14.14	3000	J
13.	UNKNOWN SILOXANE	14.19	24000	J
14.	LABORATORY ARTIFACT	14.54	4500	BJ
15.	UNKNOWN SILOXANE	14.82	21000	J
16.	UNKNOWN SILOXANE	15.42	15000	J
17.	UNKNOWN	15.97	3000	J
18.	UNKNOWN SILOXANE	16.04	12000	J
19.	UNKNOWN	16.55	7600	J
20.	UNKNOWN	17.07	6100	J
21.	BENZO FLUORANTHENE	17.19	18000	J
22.	UNKNOWN	17.55	6100	J
23.	UNKNOWN	20.60	4500	J
24.	UNKNOWN	23.45	26000	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY20

Lab Name: COMPUCHEM.RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477491

Sample wt/vol: 30.40(g/ml)G

Lab File ID:

% Moisture: 35 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/07/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.7

Sulfur Cleanup: (Y/N) N

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

319-84-6-----	alpha-BHC	2.6	U
319-85-7-----	beta-BHC	1.3	JP
319-86-8-----	delta-BHC	2.6	U
58-89-9-----	gamma-BHC (Lindane)	2.6	U
76-44-8-----	Heptachlor	1.8	JPB
309-00-2-----	Aldrin	5.6	P
1024-57-3-----	Heptachlor epoxide	2.3	JP
959-98-8-----	Endosulfan I	2.6	U
60-57-1-----	Dieldrin	14	P
72-55-9-----	4,4'-DDE	11	
72-20-8-----	Endrin	34	P
33213-65-9-----	Endosulfan II	20	P
72-54-8-----	4,4'-DDD	5.0	U
1031-07-8-----	Endosulfan sulfate	5.0	U
50-29-3-----	4,4'-DDT	5.0	U
72-43-5-----	Methoxychlor	120	PB
53494-70-5-----	Endrin ketone	5.0	U
7421-93-4-----	Endrin aldehyde	17	P
5103-71-9-----	alpha-Chlordane	2.6	U
5103-74-2-----	gamma-Chlordane	2.8	P
8001-35-2-----	Toxaphene	260	U
12674-11-2-----	Aroclor-1016	50	U
11104-28-2-----	Aroclor-1221	100	U
11141-16-5-----	Aroclor-1232	50	U
53469-21-9-----	Aroclor-1242	50	U
12672-29-6-----	Aroclor-1248	50	U
11097-69-1-----	Aroclor-1254	50	U
11096-82-5-----	Aroclor-1260	50	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY21

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477492

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077492A18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 35 Date Analyzed: 01/31/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	15	U
74-83-9-----	Bromomethane	15	U
75-01-4-----	Vinyl Chloride	15	U
75-00-3-----	Chloroethane	15	U
75-09-2-----	Methylene Chloride	51	B
67-64-1-----	Acetone	86	B
75-15-0-----	Carbon Disulfide	15	U
75-35-4-----	1,1-Dichloroethene	15	U
75-34-3-----	1,1-Dichloroethane	15	U
540-59-0-----	1,2-Dichloroethene (total)	15	U
67-66-3-----	Chloroform	15	U
107-06-2-----	1,2-Dichloroethane	15	U
78-93-3-----	2-Butanone	15	U
71-55-6-----	1,1,1-Trichloroethane	15	U
56-23-5-----	Carbon Tetrachloride	15	U
75-27-4-----	Bromodichloromethane	15	U
78-87-5-----	1,2-Dichloropropane	15	U
10061-01-5-----	cis-1,3-Dichloropropene	15	U
79-01-6-----	Trichloroethene	15	U
124-48-1-----	Dibromochloromethane	15	U
79-00-5-----	1,1,2-Trichloroethane	15	U
71-43-2-----	Benzene	15	U
10061-02-6-----	Trans-1,3-Dichloropropene	15	U
75-25-2-----	Bromoform	15	U
108-10-1-----	4-Methyl-2-Pentanone	15	U
591-78-6-----	2-Hexanone	15	U
127-18-4-----	Tetrachloroethene	15	U
79-34-5-----	1,1,2,2-Tetrachloroethane	15	U
108-88-3-----	Toluene	15	U
108-90-7-----	Chlorobenzene	15	U
100-41-4-----	Ethylbenzene	15	U
100-42-5-----	Styrene	15	U
1330-20-7-----	Xylene (total)	15	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

189

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY21

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477492

Sample wt/vol: 30.4 (g/mL) G Lab File ID: G2D77492A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 35 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 5.0

GPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND		
108-95-2-----	Phenol	2500	U
111-44-4-----	bis(2-Chloroethyl) Ether	2500	U
95-57-8-----	2-Chlorophenol	2500	U
541-73-1-----	1,3-Dichlorobenzene	2500	U
106-46-7-----	1,4-Dichlorobenzene	2500	U
95-50-1-----	1,2-Dichlorobenzene	2500	U
95-48-7-----	2-Methylphenol	2500	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	2500	U
106-44-5-----	4-Methylphenol	2500	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	2500	U
67-72-1-----	Hexachloroethane	2500	U
98-95-3-----	Nitrobenzene	2500	U
78-59-1-----	Isophorone	2500	U
88-75-5-----	2-Nitrophenol	2500	U
105-67-9-----	2,4-Dimethylphenol	2500	U
111-91-1-----	bis(2-Chloroethoxy) Methane	2500	U
120-83-2-----	2,4-Dichlorophenol	2500	U
120-82-1-----	1,2,4-Trichlorobenzene	2500	U
91-20-3-----	Naphthalene	2500	U
106-47-8-----	4-Chloroaniline	2500	U
87-68-3-----	Hexachlorobutadiene	2500	U
59-50-7-----	4-Chloro-3-Methylphenol	2500	U
91-57-6-----	2-Methylnaphthalene	2500	U
77-47-4-----	Hexachlorocyclopentadiene	2500	U
88-06-2-----	2,4,6-Trichlorophenol	2500	U
95-95-4-----	2,4,5-Trichlorophenol	6100	U
91-58-7-----	2-Chloronaphthalene	2500	U
88-74-4-----	2-Nitroaniline	6100	U
131-11-3-----	Dimethyl Phthalate	2500	U
208-96-8-----	Acenaphthylene	2500	U
606-20-2-----	2,6-Dinitrotoluene	2500	U
99-09-2-----	3-Nitroaniline	6100	U
83-32-9-----	Acenaphthene	350	J

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1258

ORIGINAL

For

1C

## SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY21

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
 Matrix: (soil/water) SOIL Lab Sample ID: 477492  
 Sample wt/vol: 30.4 (g/mL) G Lab File ID: G2D77492A21  
 Level: (low/med) LOW Date Received: 01/29/92  
 % Moisture: 35 decanted: (Y/N) N Date Extracted: 02/04/92  
 Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92  
 Injection Volume: 2.0 (uL) Dilution Factor: 5.0  
 GPC Cleanup: (Y/N) Y pH: 6.6

## CONCENTRATION UNITS:

(ug/L or ug/Kg) UG/KG

Q

CAS NO.	COMPOUND		
51-28-5-----	2,4-Dinitrophenol	6100	U
100-02-7-----	4-Nitrophenol	6100	U
132-64-9-----	Dibenzofuran	2500	U
121-14-2-----	2,4-Dinitrotoluene	2500	U
84-66-2-----	Diethylphthalate	2500	U
7005-72-3-----	4-Chlorophenyl-phenylether	2500	U
86-73-7-----	Fluorene	2500	U
100-01-6-----	4-Nitroaniline	6100	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	6100	U
86-30-6-----	N-Nitrosodiphenylamine (1)	2500	U
101-55-3-----	4-Bromophenyl-phenylether	2500	U
118-74-1-----	Hexachlorobenzene	2500	U
87-86-5-----	Pentachlorophenol	6100	U
85-01-8-----	Phenanthrene	3500	
120-12-7-----	Anthracene	920	J
86-74-8-----	Carbazole	710	J
84-74-2-----	Di-n-Butylphthalate	2500	U
206-44-0-----	Fluoranthene	8800	
129-00-0-----	Pyrene	3500	
85-68-7-----	Butylbenzylphthalate	470	J
91-94-1-----	3,3'-Dichlorobenzidine	2500	U
56-55-3-----	Benzo(a)Anthracene	4500	
218-01-9-----	Chrysene	3800	
117-81-7-----	bis(2-Ethylhexyl)Phthalate	2500	U
117-84-0-----	Di-n-Octyl Phthalate	2500	U
205-99-2-----	Benzo(b)Fluoranthene	11000	X
207-08-9-----	Benzo(k)Fluoranthene	11000	X
50-32-8-----	Benzo(a)Pyrene	6900	
193-39-5-----	Indeno(1,2,3-cd)Pyrene	3800	
53-70-3-----	Dibenz(a,h)Anthracene	1700	J
191-24-2-----	Benzo(g,h,i)Perylene	3500	

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1259

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY21

Lab Name: COMPUCHEM.RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 17744

SAS No.: 6579HQ

SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477492

Sample wt/vol: 30.4 (g/mL) G

Lab File ID: G2D77492A21

Level: (low/med) LOW

Date Received: 01/29/92

% Moisture: 35 decanted: (Y/N) N

Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL)

Dilution Factor: 5.0

GPC Cleanup: (Y/N) Y

pH: 6.6

Number TICs found: 21

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	ALDOL	4.80	1500	ABJ
2.	UNKNOWN PAH	12.45	500	J
3.	UNKNOWN	16.45	500	J
4.	UNKNOWN	16.67	760	J
5.	UNKNOWN	17.04	1300	J
6.	UNKNOWN	17.20	1000	J
7.	UNKNOWN	17.25	1000	J
8.	UNKNOWN PAH	17.34	5000	J
9.	UNKNOWN	17.90	760	J
10.	UNKNOWN	17.95	1000	J
11.	UNKNOWN	18.07	1000	J
12.	UNKNOWN	18.14	1500	J
13.	UNKNOWN	18.22	500	J
14.	UNKNOWN	18.75	1500	J
15.	UNKNOWN	18.80	1000	J
16.	UNKNOWN	19.45	500	J
17.	UNKNOWN	19.50	500	J
18.	UNKNOWN	19.55	1000	J
19.	UNKNOWN	23.39	1800	J
20.	UNKNOWN	23.42	1000	J
21.	UNKNOWN	23.45	760	J

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY22

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477493

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077493A18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 20 Date Analyzed: 01/31/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:            (uL) Soil Aliquot Volume:            (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND Q

74-87-3-----	Chloromethane	12	U
74-83-9-----	Bromomethane	12	U
75-01-4-----	Vinyl Chloride	12	U
75-00-3-----	Chloroethane	12	U
75-09-2-----	Methylene Chloride	56	B
67-64-1-----	Acetone	61	B
75-15-0-----	Carbon Disulfide	12	U
75-35-4-----	1,1-Dichloroethene	12	U
75-34-3-----	1,1-Dichloroethane	12	U
540-59-0-----	1,2-Dichloroethene (total)	12	U
67-66-3-----	Chloroform	12	U
107-06-2-----	1,2-Dichloroethane	12	U
78-93-3-----	2-Butanone	12	U
71-55-6-----	1,1,1-Trichloroethane	12	U
56-23-5-----	Carbon Tetrachloride	12	U
75-27-4-----	Bromodichloromethane	12	U
78-87-5-----	1,2-Dichloropropane	12	U
10061-01-5-----	cis-1,3-Dichloropropene	12	U
79-01-6-----	Trichloroethene	12	U
124-48-1-----	Dibromochloromethane	12	U
79-00-5-----	1,1,2-Trichloroethane	12	U
71-43-2-----	Benzene	12	U
10061-02-6-----	Trans-1,3-Dichloropropene	12	U
75-25-2-----	Bromoform	12	U
108-10-1-----	4-Methyl-2-Pentanone	12	U
591-78-6-----	2-Hexanone	12	U
127-18-4-----	Tetrachloroethene	12	U
79-34-5-----	1,1,2,2-Tetrachloroethane	12	U
108-88-3-----	Toluene	12	U
108-90-7-----	Chlorobenzene	12	U
100-41-4-----	Ethylbenzene	12	U
100-42-5-----	Styrene	12	U
1330-20-7-----	Xylene (total)	12	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

201



1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY22

Lab Name: COMPUCHEM.RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 17744

SAS No.: 6579HQ

SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477493

Sample wt/vol: 30.5 (g/mL) G

Lab File ID: GRJ77493A21

Level: (low/med) LOW

Date Received: 01/29/92

% Moisture: 20 decanted: (Y/N) N

Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 02/12/92

Injection Volume: 2.0(uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 6.4

CONCENTRATION UNITS:

CAS NO.

COMPOUND

(ug/L or ug/Kg) UG/KG

Q

108-95-2-----	Phenol	410	U
111-44-4-----	bis(2-Chloroethyl) Ether	410	U
95-57-8-----	2-Chlorophenol	410	U
541-73-1-----	1,3-Dichlorobenzene	410	U
106-46-7-----	1,4-Dichlorobenzene	410	U
95-50-1-----	1,2-Dichlorobenzene	410	U
95-48-7-----	2-Methylphenol	410	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	410	U
106-44-5-----	4-Methylphenol	410	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	410	U
67-72-1-----	Hexachloroethane	410	U
98-95-3-----	Nitrobenzene	410	U
78-59-1-----	Isophorone	410	U
88-75-5-----	2-Nitrophenol	410	U
105-67-9-----	2,4-Dimethylphenol	410	U
111-91-1-----	bis(2-Chloroethoxy) Methane	410	U
120-83-2-----	2,4-Dichlorophenol	410	U
120-82-1-----	1,2,4-Trichlorobenzene	410	U
91-20-3-----	Naphthalene	59	J
106-47-8-----	4-Chloroaniline	410	U
87-68-3-----	Hexachlorobutadiene	410	U
59-50-7-----	4-Chloro-3-Methylphenol	410	U
91-57-6-----	2-Methylnaphthalene	75	J
77-47-4-----	Hexachlorocyclopentadiene	410	U
88-06-2-----	2,4,6-Trichlorophenol	410	U
95-95-4-----	2,4,5-Trichlorophenol	980	U
91-58-7-----	2-Chloronaphthalene	410	U
88-74-4-----	2-Nitroaniline	980	U
131-11-3-----	Dimethyl Phthalate	410	U
208-96-8-----	Acenaphthylene	410	U
606-20-2-----	2,6-Dinitrotoluene	410	U
99-09-2-----	3-Nitroaniline	980	U
83-32-9-----	Acenaphthene	410	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1324

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY22

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477493

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GRJ77493A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 20 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.4

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND		Q
51-28-5-----	2,4-Dinitrophenol	980	U
100-02-7-----	4-Nitrophenol	980	U
132-64-9-----	Dibenzofuran	410	U
121-14-2-----	2,4-Dinitrotoluene	410	U
84-66-2-----	Diethylphthalate	410	U
7005-72-3-----	4-Chlorophenyl-phenylether	410	U
86-73-7-----	Fluorene	410	U
100-01-6-----	4-Nitroaniline	980	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	980	U
86-30-6-----	N-Nitrosodiphenylamine (1)	410	U
101-55-3-----	4-Bromophenyl-phenylether	410	U
118-74-1-----	Hexachlorobenzene	410	U
87-86-5-----	Pentachlorophenol	980	U
85-01-8-----	Phenanthrene	370	J
120-12-7-----	Anthracene	55	J
86-74-8-----	Carbazole	68	J
84-74-2-----	Di-n-Butylphthalate	410	U
206-44-0-----	Fluoranthene	1100	
129-00-0-----	Pyrene	530	
85-68-7-----	Butylbenzylphthalate	410	U
91-94-1-----	3,3'-Dichlorobenzidine	410	U
56-55-3-----	Benzo(a)Anthracene	410	
218-01-9-----	Chrysene	440	
117-81-7-----	bis(2-Ethylhexyl)Phthalate	410	U
117-84-0-----	Di-n-Octyl Phthalate	410	U
205-99-2-----	Benzo(b)Fluoranthene	1200	X
207-08-9-----	Benzo(k)Fluoranthene	1200	X
50-32-8-----	Benzo(a)Pyrene	380	J
193-39-5-----	Indeno(1,2,3-cd)Pyrene	450	
53-70-3-----	Dibenz(a,h)Anthracene	98	J
191-24-2-----	Benzo(g,h,i)Perylene	350	J

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1325

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY22

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477493

Sample wt/vol: 30.5 (g/mL) G Lab File ID: GRJ77493A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 20 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.4

Number TICs found: 21

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.22	410	J
2.	UNKNOWN ALKENE	4.35	570	J
3.	UNKNOWN ALKENE	4.50	4500	J
4.	UNKNOWN	4.62	820	J
5.	UNKNOWN	4.70	2000	J
6.	ALDOL	4.83	980	ABJ
7.	UNKNOWN	4.98	410	J
8.	UNKNOWN	5.07	570	J
9.	UNKNOWN	5.30	2600	J
10.	UNKNOWN	5.95	1400	J
11.	UNKNOWN	8.27	160	J
12.	UNKNOWN	16.10	160	J
13.	UNKNOWN	16.60	200	J
14.	UNKNOWN	17.22	570	J
15.	UNKNOWN	17.25	370	J
16.	UNKNOWN PAH	17.37	530	J
17.	UNKNOWN	18.32	330	J
18.	UNKNOWN	18.84	410	J
19.	UNKNOWN	18.89	570	J
20.	UNKNOWN	21.54	700	J
21.	UNKNOWN SILOXANE	23.45	250	J

ORIGINAL 1B  
(Red) SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY22RE

Lab Name: COMPUCHEM.RTP Contract: 68D10083  
 Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
 Matrix: (soil/water) SOIL Lab Sample ID: 477493  
 Sample wt/vol: 30.3 (g/mL) G Lab File ID: GRJ77493A05  
 Level: (low/med) LOW Date Received: 01/29/92  
 % Moisture: 20 decanted: (Y/N) N Date Extracted: 02/14/92  
 Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/24/92  
 Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
 GPC Cleanup: (Y/N) Y pH: 6.4

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

108-95-2-----	Phenol	410	U
111-44-4-----	bis(2-Chloroethyl) Ether	410	U
95-57-8-----	2-Chlorophenol	410	U
541-73-1-----	1,3-Dichlorobenzene	410	U
106-46-7-----	1,4-Dichlorobenzene	410	U
95-50-1-----	1,2-Dichlorobenzene	410	U
95-48-7-----	2-Methylphenol	410	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	410	U
106-44-5-----	4-Methylphenol	410	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	410	U
67-72-1-----	Hexachloroethane	410	U
98-95-3-----	Nitrobenzene	410	U
78-59-1-----	Isophorone	410	U
88-75-5-----	2-Nitrophenol	410	U
105-67-9-----	2,4-Dimethylphenol	410	U
111-91-1-----	bis(2-Chloroethoxy) Methane	410	U
120-83-2-----	2,4-Dichlorophenol	410	U
120-82-1-----	1,2,4-Trichlorobenzene	410	U
91-20-3-----	Naphthalene	71	J
106-47-8-----	4-Chloroaniline	410	U
87-68-3-----	Hexachlorobutadiene	410	U
59-50-7-----	4-Chloro-3-Methylphenol	410	U
91-57-6-----	2-Methylnaphthalene	57	J
77-47-4-----	Hexachlorocyclopentadiene	410	U
88-06-2-----	2,4,6-Trichlorophenol	410	U
95-95-4-----	2,4,5-Trichlorophenol	990	U
91-58-7-----	2-Chloronaphthalene	410	U
88-74-4-----	2-Nitroaniline	990	U
131-11-3-----	Dimethyl Phthalate	410	U
208-96-8-----	Acenaphthylene	410	U
606-20-2-----	2,6-Dinitrotoluene	410	U
99-09-2-----	3-Nitroaniline	990	U
83-32-9-----	Acenaphthene	150	J

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1389

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY22RE

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477493

Sample wt/vol: 30.3 (g/mL) G Lab File ID: GRJ77493A05

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 20 decanted: (Y/N) N Date Extracted: 02/14/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/24/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

PC Cleanup: (Y/N) Y pH: 6.4

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO. COMPOUND Q

51-28-5-----	2,4-Dinitrophenol	990	U
100-02-7-----	4-Nitrophenol	990	U
132-64-9-----	Dibenzofuran	61	J
121-14-2-----	2,4-Dinitrotoluene	410	U
84-66-2-----	Diethylphthalate	410	U
7005-72-3-----	4-Chlorophenyl-phenylether	410	U
86-73-7-----	Fluorene	89	J
100-01-6-----	4-Nitroaniline	990	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	990	U
86-30-6-----	N-Nitrosodiphenylamine (1)	410	U
101-55-3-----	4-Bromophenyl-phenylether	410	U
118-74-1-----	Hexachlorobenzene	410	U
87-86-5-----	Pentachlorophenol	990	U
85-01-8-----	Phenanthrene	900	
120-12-7-----	Anthracene	150	J
86-74-8-----	Carbazole	120	J
84-74-2-----	Di-n-Butylphthalate	410	U
206-44-0-----	Fluoranthene	1500	
129-00-0-----	Pyrene	1300	
85-68-7-----	Butylbenzylphthalate	410	U
91-94-1-----	3,3'-Dichlorobenzidine	410	U
56-55-3-----	Benzo(a) Anthracene	730	
218-01-9-----	Chrysene	880	
117-81-7-----	bis(2-Ethylhexyl) Phthalate	66	J
117-84-0-----	Di-n-Octyl Phthalate	410	U
205-99-2-----	Benzo(b) Fluoranthene	1600	
207-08-9-----	Benzo(k) Fluoranthene	570	
50-32-8-----	Benzo(a) Pyrene	550	
193-39-5-----	Indeno(1,2,3-cd) Pyrene	970	
53-70-3-----	Dibenz(a,h) Anthracene	180	J
191-24-2-----	Benzo(g,h,i) Perylene	810	

(1) - Cannot be separated from Diphenylamine

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY22RE

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HO SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477493

Sample wt/vol: 30.3 (g/mL) G Lab File ID: GRJ77493A05

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 20 decanted: (Y/N) N Date Extracted: 02/14/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/24/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.4

Number TICs found: 23 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	SOLVENT CONTAMINANT	5.37	210	BJ
2.	UNKNOWN	5.67	450	J
3.	UNKNOWN ALKENE	5.72	330	J
4.	UNKNOWN ALKENE	5.85	2800	J
5.	UNKNOWN	5.88	370	J
6.	TETRACHLOROETHANE	6.02	1000	J
7.	UNKNOWN	6.08	540	J
8.	UNKNOWN	6.17	290	J
9.	ALDOL	6.22	660	ABJ
10.	ALDOL	6.33	370	ABJ
11.	UNKNOWN	6.42	740	J
12.	UNKNOWN	6.62	2000	J
13.	UNKNOWN SILOXANE	6.95	160	J
14.	UNKNOWN	7.30	1000	J
15.	UNKNOWN	8.18	160	J
16.	UNKNOWN	8.54	160	J
17.	UNKNOWN PAH	13.70	160	J
18.	UNKNOWN PNA	13.97	82	J
19.	UNKNOWN PAH	15.25	120	J
20.	UNKNOWN PAH	15.35	120	J
21.	UNKNOWN	17.67	210	J
22.	UNKNOWN	19.74	1100	J
23.	UNKNOWN PAH	20.45	910	J

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY22

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477493

Sample wt/vol: 30.30(g/ml)G

Lab File ID:

% Moisture: 20 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/08/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.4

Sulfur Cleanup: (Y/N) N

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

319-84-6-----	alpha-BHC	2.1	U
319-85-7-----	beta-BHC	2.1	U
319-86-8-----	delta-BHC	2.1	U
58-89-9-----	gamma-BHC (Lindane)	2.1	U
76-44-8-----	Heptachlor	2.1	U
309-00-2-----	Aldrin	2.1	U
1024-57-3-----	Heptachlor epoxide	2.1	U
959-98-8-----	Endosulfan I	2.1	U
60-57-1-----	Dieldrin	6.2	P
72-55-9-----	4,4'-DDE	1.6	J
72-20-8-----	Endrin	4.1	U
33213-65-9-----	Endosulfan II	4.1	U
72-54-8-----	4,4'-DDD	4.1	U
1031-07-8-----	Endosulfan sulfate	4.1	U
50-29-3-----	4,4'-DDT	2.3	JP
72-43-5-----	Methoxychlor	26	PB
53494-70-5-----	Endrin ketone	0.66	JP
7421-93-4-----	Endrin aldehyde	4.1	U
5103-71-9-----	alpha-Chlordane	3.5	P
5103-74-2-----	gamma-Chlordane	2.1	P
8001-35-2-----	Toxaphene	210	U
12674-11-2-----	Aroclor-1016	41	U
11104-28-2-----	Aroclor-1221	83	U
11141-16-5-----	Aroclor-1232	41	U
53469-21-9-----	Aroclor-1242	41	U
12672-29-6-----	Aroclor-1248	41	U
11097-69-1-----	Aroclor-1254	41	U
11096-82-5-----	Aroclor-1260	41	U

ORIGINAL  
(Red)

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY23

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11  
Matrix: (soil/water) SOIL Lab Sample ID: 477494  
Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077494C18  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec. 22 Date Analyzed: 02/03/92  
GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	13	U
74-83-9-----	Bromomethane	13	U
75-01-4-----	Vinyl Chloride	13	U
75-00-3-----	Chloroethane	13	U
75-09-2-----	Methylene Chloride	68	B
67-64-1-----	Acetone	37	B
75-15-0-----	Carbon Disulfide	13	U
75-35-4-----	1,1-Dichloroethene	13	U
75-34-3-----	1,1-Dichloroethane	13	U
540-59-0-----	1,2-Dichloroethene (total)	13	U
67-66-3-----	Chloroform	13	U
107-06-2-----	1,2-Dichloroethane	13	U
78-93-3-----	2-Butanone	13	U
71-55-6-----	1,1,1-Trichloroethane	13	U
56-23-5-----	Carbon Tetrachloride	13	U
75-27-4-----	Bromodichloromethane	13	U
78-87-5-----	1,2-Dichloropropane	13	U
10061-01-5-----	cis-1,3-Dichloropropene	13	U
79-01-6-----	Trichloroethene	13	U
124-48-1-----	Dibromochloromethane	13	U
79-00-5-----	1,1,2-Trichloroethane	13	U
71-43-2-----	Benzene	13	U
10061-02-6-----	Trans-1,3-Dichloropropene	13	U
75-25-2-----	Bromoform	13	U
108-10-1-----	4-Methyl-2-Pentanone	13	U
591-78-6-----	2-Hexanone	13	U
127-18-4-----	Tetrachloroethene	13	U
79-34-5-----	1,1,2,2-Tetrachloroethane	13	U
108-88-3-----	Toluene	13	U
108-90-7-----	Chlorobenzene	13	U
100-41-4-----	Ethylbenzene	13	U
100-42-5-----	Styrene	13	U
1330-20-7-----	Xylene (total)	13	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

213



1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY23

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477494

Sample wt/vol: 30.5 (g/mL) G Lab File ID: G2J77494A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 22 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	Q
108-95-2-----	Phenol	420 U
111-44-4-----	bis(2-Chloroethyl) Ether	420 U
95-57-8-----	2-Chlorophenol	420 U
541-73-1-----	1,3-Dichlorobenzene	420 U
106-46-7-----	1,4-Dichlorobenzene	420 U
95-50-1-----	1,2-Dichlorobenzene	420 U
95-48-7-----	2-Methylphenol	420 U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	420 U
106-44-5-----	4-Methylphenol	420 U
621-64-7-----	N-Nitroso-Di-n-Propylamine	420 U
67-72-1-----	Hexachloroethane	420 U
98-95-3-----	Nitrobenzene	420 U
78-59-1-----	Isophorone	420 U
88-75-5-----	2-Nitrophenol	420 U
105-67-9-----	2,4-Dimethylphenol	420 U
111-91-1-----	bis(2-Chloroethoxy) Methane	420 U
120-83-2-----	2,4-Dichlorophenol	420 U
120-82-1-----	1,2,4-Trichlorobenzene	420 U
91-20-3-----	Naphthalene	420 U
106-47-8-----	4-Chloroaniline	420 U
87-68-3-----	Hexachlorobutadiene	420 U
59-50-7-----	4-Chloro-3-Methylphenol	420 U
91-57-6-----	2-Methylnaphthalene	420 U
77-47-4-----	Hexachlorocyclopentadiene	420 U
88-06-2-----	2,4,6-Trichlorophenol	420 U
95-95-4-----	2,4,5-Trichlorophenol	1000 U
91-58-7-----	2-Chloronaphthalene	420 U
88-74-4-----	2-Nitroaniline	1000 U
131-11-3-----	Dimethyl Phthalate	420 U
208-96-8-----	Acenaphthylene	420 U
606-20-2-----	2,6-Dinitrotoluene	420 U
99-09-2-----	3-Nitroaniline	1000 U
83-32-9-----	Acenaphthene	420 U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1463

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY23

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477494

Sample wt/vol: 30.5 (g/mL) G Lab File ID: G2J77494A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 22 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.

COMPOUND

Q

51-28-5-----	2,4-Dinitrophenol	1000	U
100-02-7-----	4-Nitrophenol	1000	U
132-64-9-----	Dibenzofuran	420	U
121-14-2-----	2,4-Dinitrotoluene	420	U
84-66-2-----	Diethylphthalate	54	J
7005-72-3-----	4-Chlorophenyl-phenylether	420	U
86-73-7-----	Fluorene	420	U
100-01-6-----	4-Nitroaniline	1000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1000	U
86-30-6-----	N-Nitrosodiphenylamine (1)	420	U
101-55-3-----	4-Bromophenyl-phenylether	420	U
118-74-1-----	Hexachlorobenzene	420	U
87-86-5-----	Pentachlorophenol	1000	U
85-01-8-----	Phenanthrene	420	U
120-12-7-----	Anthracene	420	U
86-74-8-----	Carbazole	420	U
84-74-2-----	Di-n-Butylphthalate	420	U
206-44-0-----	Fluoranthene	79	J
129-00-0-----	Pyrene	63	J
85-68-7-----	Butylbenzylphthalate	420	U
91-94-1-----	3,3'-Dichlorobenzidine	420	U
56-55-3-----	Benzo(a)Anthracene	420	U
218-01-9-----	Chrysene	44	J
117-81-7-----	bis(2-Ethylhexyl)Phthalate	420	U
117-84-0-----	Di-n-Octyl Phthalate	420	U
205-99-2-----	Benzo(b)Fluoranthene	84	JX
207-08-9-----	Benzo(k)Fluoranthene	84	JX
50-32-8-----	Benzo(a)Pyrene	420	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	420	U
53-70-3-----	Dibenz(a,h)Anthracene	420	U
191-24-2-----	Benzo(g,h,i)Perylene	420	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1464

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY23

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477494

Sample wt/vol: 30.5 (g/mL) G Lab File ID: G2J77494A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 22 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

PC Cleanup: (Y/N) Y pH: 6.6

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

Number TICs found: 29

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	4.40	290	J
2.	UNKNOWN	4.62	590	J
3.	ALDOL	4.77	500	ABJ
4.	ALDOL	4.88	210	AJ
5.	BENZALDEHYDE + UNKNOWN	4.98	84	J
6.	ALDOL	5.20	970	AJ
7.	ALDOL	5.38	130	AJ
8.	UNKNOWN	5.85	1300	J
9.	BLANK CONTAMINANT	9.12	130	BJ
10.	BLANK CONTAMINANT	10.24	210	BJ
11.	BLANK CONTAMINANT	11.19	340	BJ
12.	UNKNOWN SILOXANE	12.05	420	J
13.	UNKNOWN	12.22	84	J
14.	UNKNOWN	12.27	210	J
15.	UNKNOWN CARBOXYLIC ACID	12.32	340	J
16.	UNKNOWN	12.59	130	J
17.	UNKNOWN SILOXANE	12.84	340	J
18.	BLANK CONTAMINANT	13.54	800	BJ
19.	UNKNOWN SILOXANE	14.19	590	J
20.	LABORATORY ARTIFACT	14.54	210	BJ
21.	UNKNOWN SILOXANE	14.82	500	J
22.	UNKNOWN	15.07	170	J
23.	UNKNOWN SILOXANE	15.42	380	J
24.	UNKNOWN SILOXANE	16.02	550	J
25.	UNKNOWN	16.70	630	J
26.	UNKNOWN	17.05	930	J
27.	UNKNOWN	17.15	420	J
28.	UNKNOWN	18.65	2200	J
29.	UNKNOWN	21.25	880	J

FORM I SV-TIC

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1465

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY23

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477494

Sample wt/vol: 30.30(g/ml)G

Lab File ID:

% Moisture: 22 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/08/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.6

Sulfur Cleanup: (Y/N) N

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
319-84-6-----	alpha-BHC	2.2	U
319-85-7-----	beta-BHC	0.87	JP
319-86-8-----	delta-BHC	2.2	U
58-89-9-----	gamma-BHC (Lindane)	2.2	U
76-44-8-----	Heptachlor	0.23	JPB
309-00-2-----	Aldrin	2.2	U
1024-57-3-----	Heptachlor epoxide	2.2	U
959-98-8-----	Endosulfan I	2.2	U
60-57-1-----	Dieldrin	4.2	U
72-55-9-----	4,4'-DDE	4.2	U
72-20-8-----	Endrin	1.4	J
33213-65-9-----	Endosulfan II	0.41	JP
72-54-8-----	4,4'-DDD	4.2	U
1031-07-8-----	Endosulfan sulfate	4.2	U
50-29-3-----	4,4'-DDT	4.2	U
72-43-5-----	Methoxychlor	1.4	JPB
53494-70-5-----	Endrin ketone	4.2	U
7421-93-4-----	Endrin aldehyde	0.68	JP
5103-71-9-----	alpha-Chlordane	2.2	U
5103-74-2-----	gamma-Chlordane	2.2	U
8001-35-2-----	Toxaphene	220	U
12674-11-2-----	Aroclor-1016	42	U
11104-28-2-----	Aroclor-1221	85	U
11141-16-5-----	Aroclor-1232	42	U
53469-21-9-----	Aroclor-1242	42	U
12672-29-6-----	Aroclor-1248	42	U
11097-69-1-----	Aroclor-1254	42	U
11096-82-5-----	Aroclor-1260	42	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY24

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477467

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CR077467C51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/03/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

74-87-3-----	Chloromethane	10	U
74-83-9-----	Bromomethane	10	U
75-01-4-----	Vinyl Chloride	10	U
75-00-3-----	Chloroethane	10	U
75-09-2-----	Methylene Chloride	160	B
67-64-1-----	Acetone	10	U
75-15-0-----	Carbon Disulfide	10	U
75-35-4-----	1,1-Dichloroethene	10	U
75-34-3-----	1,1-Dichloroethane	10	U
540-59-0-----	1,2-Dichloroethene (total)	10	U
67-66-3-----	Chloroform	10	U
107-06-2-----	1,2-Dichloroethane	10	U
78-93-3-----	2-Butanone	10	U
71-55-6-----	1,1,1-Trichloroethane	10	U
56-23-5-----	Carbon Tetrachloride	10	U
75-27-4-----	Bromodichloromethane	10	U
78-87-5-----	1,2-Dichloropropane	10	U
10061-01-5-----	cis-1,3-Dichloropropene	10	U
79-01-6-----	Trichloroethene	10	U
124-48-1-----	Dibromochloromethane	10	U
79-00-5-----	1,1,2-Trichloroethane	10	U
71-43-2-----	Benzene	10	U
10061-02-6-----	Trans-1,3-Dichloropropene	10	U
75-25-2-----	Bromoform	10	U
108-10-1-----	4-Methyl-2-Pentanone	10	U
591-78-6-----	2-Hexanone	10	U
127-18-4-----	Tetrachloroethene	10	U
79-34-5-----	1,1,2,2-Tetrachloroethane	10	U
108-88-3-----	Toluene	10	U
108-90-7-----	Chlorobenzene	10	U
100-41-4-----	Ethylbenzene	10	U
100-42-5-----	Styrene	10	U
1330-20-7-----	Xylene (total)	10	U

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY24

Lab Name: COMPUCHEM, RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477467  
Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CR077467C51  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec.      Date Analyzed: 02/03/92  
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 0 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY24

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477467

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077467A07

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 01/31/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

108-95-2-----	Phenol	10	U
111-44-4-----	bis(2-Chloroethyl) Ether	10	U
95-57-8-----	2-Chlorophenol	10	U
541-73-1-----	1,3-Dichlorobenzene	10	U
106-46-7-----	1,4-Dichlorobenzene	10	U
95-50-1-----	1,2-Dichlorobenzene	10	U
95-48-7-----	2-Methylphenol	10	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	10	U
106-44-5-----	4-Methylphenol	10	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	10	U
67-72-1-----	Hexachloroethane	10	U
98-95-3-----	Nitrobenzene	10	U
78-59-1-----	Isophorone	10	U
88-75-5-----	2-Nitrophenol	10	U
105-67-9-----	2,4-Dimethylphenol	10	U
111-91-1-----	bis(2-Chloroethoxy) Methane	10	U
120-83-2-----	2,4-Dichlorophenol	10	U
120-82-1-----	1,2,4-Trichlorobenzene	10	U
91-20-3-----	Naphthalene	10	U
106-47-8-----	4-Chloroaniline	10	U
87-68-3-----	Hexachlorobutadiene	10	U
59-50-7-----	4-Chloro-3-Methylphenol	10	U
91-57-6-----	2-Methylnaphthalene	10	U
77-47-4-----	Hexachlorocyclopentadiene	10	U
88-06-2-----	2,4,6-Trichlorophenol	10	U
95-95-4-----	2,4,5-Trichlorophenol	25	U
91-58-7-----	2-Chloronaphthalene	10	U
88-74-4-----	2-Nitroaniline	25	U
131-11-3-----	Dimethyl Phthalate	10	U
208-96-8-----	Acenaphthylene	10	U
606-20-2-----	2,6-Dinitrotoluene	10	U
99-09-2-----	3-Nitroaniline	25	U
83-32-9-----	Acenaphthene	10	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

406

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY24

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477467

Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077467A07

Level: (low/med) LOW Date Received: 01/29/92

% Moisture:        decanted: (Y/N)        Date Extracted: 01/31/92

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH:       

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

51-28-5-----	2,4-Dinitrophenol	25	U
100-02-7-----	4-Nitrophenol	25	U
132-64-9-----	Dibenzofuran	10	U
121-14-2-----	2,4-Dinitrotoluene	10	U
84-66-2-----	Diethylphthalate	10	U
7005-72-3-----	4-Chlorophenyl-phenylether	10	U
86-73-7-----	Fluorene	10	U
100-01-6-----	4-Nitroaniline	25	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	25	U
86-30-6-----	N-Nitrosodiphenylamine (1)	10	U
101-55-3-----	4-Bromophenyl-phenylether	10	U
118-74-1-----	Hexachlorobenzene	10	U
87-86-5-----	Pentachlorophenol	25	U
85-01-8-----	Phenanthrene	10	U
120-12-7-----	Anthracene	10	U
86-74-8-----	Carbazole	10	U
84-74-2-----	Di-n-Butylphthalate	10	U
206-44-0-----	Fluoranthene	10	U
129-00-0-----	Pyrene	10	U
85-68-7-----	Butylbenzylphthalate	10	U
91-94-1-----	3,3'-Dichlorobenzidine	10	U
56-55-3-----	Benzo(a)Anthracene	10	U
218-01-9-----	Chrysene	10	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	10	U
117-84-0-----	Di-n-Octyl Phthalate	10	U
205-99-2-----	Benzo(b) Fluoranthene	10	U
207-08-9-----	Benzo(k) Fluoranthene	10	U
50-32-8-----	Benzo(a) Pyrene	10	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	10	U
53-70-3-----	Dibenz(a,h) Anthracene	10	U
191-24-2-----	Benzo(g,h,i) Perylene	10	U

(1) - Cannot be separated from Diphenylamine



1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY24

Lab Name: COMPUCHEM RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477467  
Sample wt/vol: 1000 (g/mL) ML Lab File ID: GH077467A07  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture:        decanted: (Y/N)        Date Extracted: 01/31/92  
Concentrated Extract Volume: 1000 (uL) Date Analyzed: 02/04/92  
Injection Volume: 2.0 (uL) Dilution Factor: 1.0  
PC Cleanup: (Y/N) N pH:       

Number TICs found: 0

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY24

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02

Matrix: (soil/water) WATER

Lab Sample ID: 477467

Sample wt/vol: 1000(g/ml)ML

Lab File ID:

% Moisture: decanted: (Y/N)

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SEPF

Date Extracted: 01/30/92

Concentrated Extract Volume: 10000(uL)

Date Analyzed: 02/04/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) N

pH:

Sulfur Cleanup: (Y/N) N

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

319-84-6-----	alpha-BHC	0.050	U
319-85-7-----	beta-BHC	0.050	U
319-86-8-----	delta-BHC	0.050	U
58-89-9-----	gamma-BHC (Lindane)	0.050	U
76-44-8-----	Heptachlor	0.050	U
309-00-2-----	Aldrin	0.050	U
1024-57-3-----	Heptachlor epoxide	0.050	U
959-98-8-----	Endosulfan I	0.050	U
60-57-1-----	Dieldrin	0.10	U
72-55-9-----	4,4'-DDE	0.10	U
72-20-8-----	Endrin	0.10	U
33213-65-9-----	Endosulfan II	0.10	U
72-54-8-----	4,4'-DDD	0.10	U
1031-07-8-----	Endosulfan sulfate	0.10	U
50-29-3-----	4,4'-DDT	0.10	U
72-43-5-----	Methoxychlor	0.50	U
53494-70-5-----	Endrin ketone	0.10	U
7421-93-4-----	Endrin aldehyde	0.10	U
5103-71-9-----	alpha-Chlordane	0.050	U
5103-74-2-----	gamma-Chlordane	0.050	U
8001-35-2-----	Toxaphene	5.0	U
12674-11-2-----	Aroclor-1016	1.0	U
11104-28-2-----	Aroclor-1221	2.0	U
11141-16-5-----	Aroclor-1232	1.0	U
53469-21-9-----	Aroclor-1242	1.0	U
12672-29-6-----	Aroclor-1248	1.0	U
11097-69-1-----	Aroclor-1254	1.0	U
11096-82-5-----	Aroclor-1260	1.0	U

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY25

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HO SDG No.: CHY02

Matrix: (soil/water) WATER Lab Sample ID: 477468

Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CR077468C51

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec.        Date Analyzed: 02/03/92

GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume:        (uL) Soil Aliquot Volume:        (uL)

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/L</u>	Q
74-87-3	-----Chloromethane	10	U
74-83-9	-----Bromomethane	10	U
75-01-4	-----Vinyl Chloride	10	U
75-00-3	-----Chloroethane	10	U
75-09-2	-----Methylene Chloride	150	B
67-64-1	-----Acetone	10	U
75-15-0	-----Carbon Disulfide	10	U
75-35-4	-----1,1-Dichloroethene	10	U
75-34-3	-----1,1-Dichloroethane	10	U
540-59-0	-----1,2-Dichloroethene (total)	10	U
67-66-3	-----Chloroform	10	U
107-06-2	-----1,2-Dichloroethane	10	U
78-93-3	-----2-Butanone	10	U
71-55-6	-----1,1,1-Trichloroethane	10	U
56-23-5	-----Carbon Tetrachloride	10	U
75-27-4	-----Bromodichloromethane	10	U
78-87-5	-----1,2-Dichloropropane	10	U
10061-01-5	-----cis-1,3-Dichloropropene	10	U
79-01-6	-----Trichloroethene	10	U
124-48-1	-----Dibromochloromethane	10	U
79-00-5	-----1,1,2-Trichloroethane	10	U
71-43-2	-----Benzene	10	U
10061-02-6	-----Trans-1,3-Dichloropropene	10	U
75-25-2	-----Bromoform	10	U
108-10-1	-----4-Methyl-2-Pentanone	10	U
591-78-6	-----2-Hexanone	10	U
127-18-4	-----Tetrachloroethene	10	U
79-34-5	-----1,1,2,2-Tetrachloroethane	10	U
108-88-3	-----Toluene	10	U
108-90-7	-----Chlorobenzene	10	U
100-41-4	-----Ethylbenzene	10	U
100-42-5	-----Styrene	10	U
1330-20-7	-----Xylene (total)	10	U

FORM I VOA

3/90

SAMPLE DATA PACKAGE

17744 CHY 02

135

1E  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY25

Lab Name: COMPUCHEM RTP Contract: 68D10083  
Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY02  
Matrix: (soil/water) WATER Lab Sample ID: 477468  
Sample wt/vol: 5.0 (g/mL) ML Lab File ID: CR077468C51  
Level: (low/med) LOW Date Received: 01/29/92  
% Moisture: not dec.      Date Analyzed: 02/03/92  
GC Column: DB-624 ID: 0.530 (mm) Dilution Factor: 1.0  
Soil Extract Volume:      (uL) Soil Aliquot Volume:      (uL)

Number TICs found: 0 CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

1A  
VOLATILE ORGANICS ANALYSIS DATA SHEET

ORIGINAL  
EPA SAMPLE NO. 68D10083

CHY26

Lab Name: COMPUCHEM, RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477495

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH077495C18

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: not dec. 32 Date Analyzed: 02/03/92

GC Column: DB624 ID: 0.530 (mm) Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG Q

74-87-3-----	Chloromethane	15	U
74-83-9-----	Bromomethane	15	U
75-01-4-----	Vinyl Chloride	15	U
75-00-3-----	Chloroethane	15	U
75-09-2-----	Methylene Chloride	83	B
67-64-1-----	Acetone	73	B
75-15-0-----	Carbon Disulfide	15	U
75-35-4-----	1,1-Dichloroethene	15	U
75-34-3-----	1,1-Dichloroethane	15	U
540-59-0-----	1,2-Dichloroethene (total)	15	U
67-66-3-----	Chloroform	15	U
107-06-2-----	1,2-Dichloroethane	15	U
78-93-3-----	Butanone	15	U
71-55-6-----	1,1,1-Trichloroethane	15	U
56-23-5-----	Carbon Tetrachloride	15	U
75-27-4-----	Bromodichloromethane	15	U
78-87-5-----	1,2-Dichloropropane	15	U
10061-01-5-----	cis-1,3-Dichloropropene	15	U
79-01-6-----	Trichloroethene	15	U
124-48-1-----	Dibromochloromethane	15	U
79-00-5-----	1,1,2-Trichloroethane	15	U
71-43-2-----	Benzene	15	U
10061-02-6-----	Trans-1,3-Dichloropropene	15	U
75-25-2-----	Bromoform	15	U
108-10-1-----	4-Methyl-2-Pentanone	15	U
591-78-6-----	2-Hexanone	15	U
127-18-4-----	Tetrachloroethene	15	U
79-34-5-----	1,1,2,2-Tetrachloroethane	15	U
108-88-3-----	Toluene	15	U
108-90-7-----	Chlorobenzene	15	U
100-41-4-----	Ethylbenzene	15	U
100-42-5-----	Styrene	15	U
1330-20-7-----	Xylene (total)	15	U

1B  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY26

Lab Name: COMPUCHEM RTP

Contract: 68D10083

Lab Code: COMPU

Case No.: 17744

SAS No.: 6579HQ

SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477495

Sample wt/vol: 30.4 (g/mL) G

Lab File ID: GRJ77495A21

Level: (low/med) LOW

Date Received: 01/29/92

% Moisture: 32 decanted: (Y/N) N

Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL)

Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y

pH: 6.3

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NO.

COMPOUND

Q

108-95-2-----	Phenol	480	U
111-44-4-----	bis(2-Chloroethyl) Ether	480	U
95-57-8-----	2-Chlorophenol	480	U
541-73-1-----	1,3-Dichlorobenzene	480	U
106-46-7-----	1,4-Dichlorobenzene	480	U
95-50-1-----	1,2-Dichlorobenzene	480	U
95-48-7-----	2-Methylphenol	480	U
108-60-1-----	2,2'-Oxybis(1-Chloropropane)	480	U
106-44-5-----	4-Methylphenol	480	U
621-64-7-----	N-Nitroso-Di-n-Propylamine	480	U
67-72-1-----	Hexachloroethane	480	U
98-95-3-----	Nitrobenzene	480	U
78-59-1-----	Isophorone	480	U
88-75-5-----	2-Nitrophenol	480	U
105-67-9-----	2,4-Dimethylphenol	480	U
111-91-1-----	bis(2-Chloroethoxy)Methane	480	U
120-83-2-----	2,4-Dichlorophenol	480	U
120-82-1-----	1,2,4-Trichlorobenzene	480	U
91-20-3-----	Naphthalene	480	U
106-47-8-----	4-Chloroaniline	480	U
87-68-3-----	Hexachlorobutadiene	480	U
59-50-7-----	4-Chloro-3-Methylphenol	480	U
91-57-6-----	2-Methylnaphthalene	480	U
77-47-4-----	Hexachlorocyclopentadiene	480	U
88-06-2-----	2,4,6-Trichlorophenol	480	U
95-95-4-----	2,4,5-Trichlorophenol	1200	U
91-58-7-----	2-Chloronaphthalene	480	U
88-74-4-----	2-Nitroaniline	1200	U
131-11-3-----	Dimethyl Phthalate	480	U
208-96-8-----	Acenaphthylene	480	U
606-20-2-----	2,6-Dinitrotoluene	480	U
99-09-2-----	3-Nitroaniline	1200	U
83-32-9-----	Acenaphthene	480	U

FORM I SV-1

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1520

1C  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY26

Lab Name: COMPUCHEM.RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477495

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GRJ77495A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 32 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

%PC Cleanup: (Y/N) Y pH: 6.3

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

51-28-5-----	2,4-Dinitrophenol	1200	U
100-02-7-----	4-Nitrophenol	1200	U
132-64-9-----	Dibenzofuran	480	U
121-14-2-----	2,4-Dinitrotoluene	480	U
84-66-2-----	Diethylphthalate	480	U
7005-72-3-----	4-Chlorophenyl-phenylether	480	U
86-73-7-----	Fluorene	480	U
100-01-6-----	4-Nitroaniline	1200	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1200	U
86-30-6-----	N-Nitrosodiphenylamine (1)	480	U
101-55-3-----	4-Bromophenyl-phenylether	480	U
118-74-1-----	Hexachlorobenzene	480	U
87-86-5-----	Pentachlorophenol	1200	U
85-01-8-----	Phenanthrene	77	J
120-12-7-----	Anthracene	480	U
86-74-8-----	Carbazole	480	U
84-74-2-----	Di-n-Butylphthalate	480	U
206-44-0-----	Fluoranthene	170	J
129-00-0-----	Pyrene	99	J
85-68-7-----	Butylbenzylphthalate	480	U
91-94-1-----	3,3'-Dichlorobenzidine	480	U
56-55-3-----	Benzo(a)Anthracene	72	J
218-01-9-----	Chrysene	86	J
117-81-7-----	bis(2-Ethylhexyl) Phthalate	480	U
117-84-0-----	Di-n-Octyl Phthalate	480	U
205-99-2-----	Benzo(b)Fluoranthene	170	JX
207-08-9-----	Benzo(k)Fluoranthene	170	JX
50-32-8-----	Benzo(a)Pyrene	86	J
193-39-5-----	Indeno(1,2,3-cd)Pyrene	480	U
53-70-3-----	Dibenz(a,h)Anthracene	480	U
191-24-2-----	Benzo(g,h,i)Perylene	480	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

3/90

SAMPLE DATA PACKAGE

17744 CHY 11

1521

1F  
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

CHY26

Lab Name: COMPUCHEM RTP Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HO SDG No.: CHY11

Matrix: (soil/water) SOIL Lab Sample ID: 477495

Sample wt/vol: 30.4 (g/mL) G Lab File ID: GRJ77495A21

Level: (low/med) LOW Date Received: 01/29/92

% Moisture: 32 decanted: (Y/N) N Date Extracted: 02/04/92

Concentrated Extract Volume: 500.0 (uL) Date Analyzed: 02/12/92

Injection Volume: 2.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) Y pH: 6.3

Number TICs found: 23

CONCENTRATION UNITS:  
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN ALKENE	4.30	190	J
2.	UNKNOWN	4.43	1500	J
3.	UNKNOWN	4.65	1100	J
4.	ALDOL	4.80	1400	ABJ
5.	ALDOL	4.92	390	AJ
6.	UNKNOWN	5.02	240	J
7.	UNKNOWN	5.25	1500	J
8.	UNKNOWN	5.90	2800	J
9. 57-10-3	HEXADECANOIC ACID	12.37	190	JN
10.	LABORATORY ARTIFACT	14.60	340	BJ
11.	UNKNOWN	15.12	240	J
12.	UNKNOWN	16.02	97	J
13.	UNKNOWN	16.05	240	J
14.	UNKNOWN	17.17	440	J
15.	UNKNOWN	17.24	290	J
16.	UNKNOWN	18.82	770	J
17.	UNKNOWN	20.47	3000	J
18.	UNKNOWN	20.49	1600	J
19.	UNKNOWN	21.50	820	J
20.	UNKNOWN	21.87	340	J
21.	UNKNOWN	21.94	190	J
22.	UNKNOWN	23.14	480	J
23.	UNKNOWN	23.42	190	J



1D  
PESTICIDE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

CHY26

Lab Name: COMPUCHEM, RTP

Contract: 68D10083

Lab Code: COMPU Case No.: 17744 SAS No.: 6579HQ SDG No.: CHY11

Matrix: (soil/water) SOIL

Lab Sample ID: 477495

Sample wt/vol: 30.20(g/ml)G

Lab File ID:

% Moisture: 32 decanted: (Y/N) N

Date Received: 01/29/92

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 02/03/92

Concentrated Extract Volume: 5000(uL)

Date Analyzed: 02/08/92

Injection Volume: 2.0(uL)

Dilution Factor: 1

GPC Cleanup: (Y/N) Y

pH: 6.3

Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
---------	----------	--	---

319-84-6-----	alpha-BHC	2.5	U
319-85-7-----	beta-BHC	0.28	JP
319-86-8-----	delta-BHC	2.5	U
58-89-9-----	gamma-BHC (Lindane)	0.091	JP
76-44-8-----	Heptachlor	0.19	JPB
309-00-2-----	Aldrin	0.56	J
1024-57-3-----	Heptachlor epoxide	2.5	U
959-98-8-----	Endosulfan I	2.5	U
60-57-1-----	Dieldrin	0.26	JP
72-55-9-----	4,4'-DDE	0.80	JP
72-20-8-----	Endrin	3.3	J
33213-65-9-----	Endosulfan II	0.28	JP
72-54-8-----	4,4'-DDD	4.8	U
1031-07-8-----	Endosulfan sulfate	4.8	U
50-29-3-----	4,4'-DDT	0.92	JP
72-43-5-----	Methoxychlor	25	U
53494-70-5-----	Endrin ketone	4.8	U
7421-93-4-----	Endrin aldehyde	4.8	U
5103-71-9-----	alpha-Chlordane	2.5	U
5103-74-2-----	gamma-Chlordane	2.5	U
8001-35-2-----	Toxaphene	250	U
12674-11-2-----	Aroclor-1016	48	U
11104-28-2-----	Aroclor-1221	98	U
11141-16-5-----	Aroclor-1232	48	U
53469-21-9-----	Aroclor-1242	48	U
12672-29-6-----	Aroclor-1248	48	U
11097-69-1-----	Aroclor-1254	48	U
11096-82-5-----	Aroclor-1260	48	U

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE02

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix (soil/water): WATER

Lab Sample ID: 9201301002

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum_	22.90	B	E	P
7440-36-0	Antimony_	30.00	U		P
7440-38-2	Arsenic_	10.90			F
7440-39-3	Barium_	85.50	B		P
7440-41-7	Beryllium	1.00	U		P
7440-43-9	Cadmium_	3.00	U		P
7440-70-2	Calcium_	68700.00			P
7440-47-3	Chromium_	3.00	U		P
7440-48-4	Cobalt_	4.00	U		P
7440-50-8	Copper_	5.00	B		P
7439-89-6	Iron_	9.40	B	NE	P
7439-92-1	Lead_	1.00	U		F
7439-95-4	Magnesium	15100.00			P
7439-96-5	Manganese	2.00	U		P
7439-97-6	Mercury_	0.20	U		CV
7439-02-0	Nickel_	22.00	U		P
7440-09-7	Potassium	774.00	B		P
7782-49-2	Selenium_	2.00	U	+	F
7440-22-4	Silver_	2.00	U		P
7440-23-5	Sodium_	14700.00			P
7440-28-0	Thallium_	2.00	U		F
7440-62-2	Vanadium_	4.00	U		P
7440-66-6	Zinc_	14.40	B		P
	Cyanide_	10.00	U		AS

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJEO3

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJEO2

Matrix (soil/water): WATER

Lab Sample ID: 9201301003

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	15.10	B	E	P
7440-36-0	Antimony	30.00	U		P
7440-38-2	Arsenic	8.10	B		F
7440-39-3	Barium	248.00			P
7440-41-7	Beryllium	1.00	U		P
7440-43-9	Cadmium	3.00	U		P
7440-70-2	Calcium	38700.00			P
7440-47-3	Chromium	3.00	U		P
7440-48-4	Cobalt	4.00	U		P
7440-50-8	Copper	188.00			P
7439-89-6	Iron	52.50	B	NE	P
7439-92-1	Lead	1.00	U		F
7439-95-4	Magnesium	15300.00			P
7439-96-5	Manganese	2.00	U		P
7439-97-6	Mercury	0.20	U		CV
7439-02-0	Nickel	22.00	U		P
7440-09-7	Potassium	1000.00	B		P
7782-49-2	Selenium	2.00	U	S	F
7440-22-4	Silver	2.00	U		P
7440-23-5	Sodium	21900.00			P
7440-28-0	Thallium	2.00	U		F
7440-62-2	Vanadium	4.00	U		P
7440-66-6	Zinc	63.30			P
	Cyanide	10.00	U		AS

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE04

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix (soil/water): WATER

Lab Sample ID: 9201301004

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	32.50	B	E	P
7440-36-0	Antimony	30.00	U		P
7440-38-2	Arsenic	5.80	B		F
7440-39-3	Barium	284.00			P
7440-41-7	Beryllium	1.00	U		P
7440-43-9	Cadmium	3.00	U		P
7440-70-2	Calcium	48400.00			P
7440-47-3	Chromium	3.00	U		P
7440-48-4	Cobalt	4.00	U		P
7440-50-8	Copper	68.00			P
7439-89-6	Iron	21.90	B	NE	P
7439-92-1	Lead	1.00	U		F
7439-95-4	Magnesium	16800.00			P
7439-96-5	Manganese	2.70	B		P
7439-97-6	Mercury	0.20	U		CV
7439-02-0	Nickel	22.00	U		P
7440-09-7	Potassium	1300.00	B		P
7782-49-2	Selenium	2.00	U	+	F
7440-22-4	Silver	2.00	U		P
7440-23-5	Sodium	17700.00			P
7440-28-0	Thallium	2.00	U		F
7440-62-2	Vanadium	4.00	U		P
7440-66-6	Zinc	39.50			P
	Cyanide	10.00	U		AS

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:

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000007

## U.S. EPA - CLP

6816.1  
(10)1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE05

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix (soil/water): WATER

Lab Sample ID: 9201301005

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	134.00	B	E	P
7440-36-0	Antimony	30.00	U		P
7440-38-2	Arsenic	6.80	B		F
7440-39-3	Barium	33.00	B		P
7440-41-7	Beryllium	1.00	U		P
7440-43-9	Cadmium	3.00	U		P
7440-70-2	Calcium	38000.00			P
7440-47-3	Chromium	3.00	U		P
7440-48-4	Cobalt	4.00	U		P
7440-50-8	Copper	5.50	B		P
7439-89-6	Iron	202.00		NE	P
7439-92-1	Lead	1.00	U		F
7439-95-4	Magnesium	12400.00			P
7439-96-5	Manganese	123.00			P
7439-97-6	Mercury	0.20	U		CV
7439-02-0	Nickel	22.00	U		P
7440-09-7	Potassium	3030.00	B		P
7782-49-2	Selenium	2.00	B		F
7440-22-4	Silver	2.00	U		P
7440-23-5	Sodium	23000.00			P
7440-28-0	Thallium	2.00	U		F
7440-62-2	Vanadium	4.00	U		P
7440-66-6	Zinc	48.50			P
	Cyanide	10.00	U		AS

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE06

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDS No.: MCJE02

Matrix (soil/water): WATER

Lab Sample ID: 9201301006

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum_	837.00		E	P
7440-36-0	Antimony_	30.00	U		P
7440-38-2	Arsenic__	7.20	B		F
7440-39-3	Barium___	55.60	B		P
7440-41-7	Beryllium	1.00	U		P
7440-43-9	Cadmium__	3.00	U		P
7440-70-2	Calcium__	38900.00			P
7440-47-3	Chromium_	3.70	B		P
7440-48-4	Cobalt___	6.70	B		P
7440-50-8	Copper___	23.80	B		P
7439-89-6	Iron_____	1380.00		NE	P
7439-92-1	Lead_____	19.00		S	F
7439-95-4	Magnesium	12700.00			P
7439-96-5	Manganese	507.00			P
7439-97-6	Mercury__	0.20	U		CV
7439-02-0	Nickel___	22.00	U		P
7440-09-7	Potassium	3300.00	B		P
7782-49-2	Selenium_	2.00	B		F
7440-22-4	Silver___	2.00	U		P
7440-23-5	Sodium___	22600.00			P
7440-28-0	Thallium_	2.00	U		F
7440-62-2	Vanadium_	4.00	U		P
7440-66-6	Zinc_____	176.00			P
	Cyanide__	11.70			AS

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:

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## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE07

b Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix (soil/water): WATER

Lab Sample ID: 9201301007

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum_	1090.00		E	P
7440-36-0	Antimony_	30.00	U		P
7440-38-2	Arsenic__	7.20	B		F
7440-39-3	Barium___	65.30	B		P
7440-41-7	Beryllium	1.00	U		P
7440-43-9	Cadmium__	3.00	U		P
7440-70-2	Calcium__	14500.00			P
7440-47-3	Chromium_	3.00	U		P
7440-48-4	Cobalt___	4.00	U		P
7440-50-8	Copper___	5.00	B		P
7439-89-6	Iron_____	930.00		NE	P
7439-92-1	Lead_____	1.10	B		F
7439-95-4	Magnesium	5790.00			P
7439-96-5	Manganese	51.70			P
7439-97-6	Mercury__	0.20	U		CV
7439-02-0	Nickel___	22.00	U		P
7440-09-7	Potassium	2270.00	B		P
7782-49-2	Selenium_	2.00	U		F
7440-22-4	Silver___	2.00	U		P
7440-23-5	Sodium___	6150.00			P
7440-28-0	Thallium_	2.00	U		F
7440-62-2	Vanadium_	4.00	U		P
7440-66-6	Zinc_____	27.70			P
	Cyanide__	10.00	U		AS

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:

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000010

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE08

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix (soil/water): WATER

Lab Sample ID: 9201301010

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	1580.00		E	P
7440-36-0	Antimony	30.00	U		P
7440-38-2	Arsenic	8.90	B		F
7440-39-3	Barium	93.90	B		P
7440-41-7	Beryllium	1.10	B		P
7440-43-9	Cadmium	3.00	U		P
7440-70-2	Calcium	44100.00			P
7440-47-3	Chromium	7.10	B		P
7440-48-4	Cobalt	29.00	B		P
7440-50-8	Copper	55.60			P
7439-89-6	Iron	3470.00		NE	P
7439-92-1	Lead	59.90			F
7439-95-4	Magnesium	14300.00			P
7439-96-5	Manganese	1210.00			P
7439-97-6	Mercury	0.20	U		CV
7439-02-0	Nickel	32.60	B		P
7440-09-7	Potassium	3770.00	B		P
7782-49-2	Selenium	2.00	U		F
7440-22-4	Silver	2.00	U		P
7440-23-5	Sodium	24100.00			P
7440-28-0	Thallium	2.00	U		F
7440-62-2	Vanadium	6.40	B		P
7440-66-6	Zinc	301.00			P
	Cyanide	10.00	U		AS

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:

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000011



## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE09

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix (soil/water): WATER

Lab Sample ID: 9201301011

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	119.00	B	E	P
7440-36-0	Antimony	30.00	U		P
7440-38-2	Arsenic	6.70	B		F
7440-39-3	Barium	30.30	B		P
7440-41-7	Beryllium	1.00	U		P
7440-43-9	Cadmium	3.00	U		P
7440-70-2	Calcium	33100.00			P
7440-47-3	Chromium	3.00	U		P
7440-48-4	Cobalt	4.00	U		P
7440-50-8	Copper	4.50	B		P
7439-89-6	Iron	132.00		NE	P
7439-92-1	Lead	1.00	U		F
7439-95-4	Magnesium	10800.00			P
7439-96-5	Manganese	113.00			P
7439-97-6	Mercury	0.20	U		CV
7439-02-0	Nickel	22.00	U		P
7440-09-7	Potassium	2780.00	B		P
7782-49-2	Selenium	2.50	B		F
7440-22-4	Silver	2.00	U		P
7440-23-5	Sodium	18800.00			P
7440-28-0	Thallium	2.00	U		F
7440-62-2	Vanadium	4.00	U		P
7440-66-6	Zinc	76.00			P
	Cyanide	10.00	U		AS

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:

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000012

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE10

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix (soil/water): WATER

Lab Sample ID: 9201301012

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

CAS No.	Analyte	Concentration	C	D	M
7429-90-5	Aluminum	266.00		E	P
7440-36-0	Antimony	30.00	U		P
7440-38-2	Arsenic	6.40	B		F
7440-39-3	Barium	33.80	B		P
7440-41-7	Beryllium	1.00	U		P
7440-43-9	Cadmium	3.00	U		P
7440-70-2	Calcium	37900.00			P
7440-47-3	Chromium	3.00	U		P
7440-48-4	Cobalt	6.40	B		P
7440-50-8	Copper	7.40	B		P
7439-89-6	Iron	547.00		NE	P
7439-92-1	Lead	2.70	B		F
7439-95-4	Magnesium	12400.00			P
7439-96-5	Manganese	379.00			P
7439-97-6	Mercury	0.20	U		CV
7439-02-0	Nickel	22.00	U		P
7440-09-7	Potassium	2620.00	B		P
7782-49-2	Selenium	3.30	B		F
7440-22-4	Silver	2.00	U		P
7440-23-5	Sodium	22300.00			P
7440-28-0	Thallium	2.00	U		F
7440-62-2	Vanadium	4.00	U		P
7440-66-6	Zinc	61.40			P
	Cyanide	10.00	U		AS

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:

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000013

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE11

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302002

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 51.2

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	6080.00			P
7440-38-0	Antimony	11.70	U		P
7440-38-2	Arsenic	7.00			F
7440-39-3	Barium	120.00			P
7440-41-7	Beryllium	1.10	B		P
7440-43-9	Cadmium	1.40	B		P
7440-70-2	Calcium	10200.00			P
7440-47-3	Chromium	31.60			P
7440-48-4	Cobalt	38.80			P
7440-50-8	Copper	74.20			P
7439-89-6	Iron	14900.00			P
7439-92-1	Lead	113.00			F
7439-95-4	Magnesium	5420.00			P
7439-96-5	Manganese	1340.00			P
7439-97-6	Mercury	0.29			CV
7439-02-0	Nickel	40.30			P
7440-09-7	Potassium	629.00	B		P
7782-49-2	Selenium	1.50	B		F
7440-22-4	Silver	0.78	U		P
7440-23-5	Sodium	203.00	B		P
7440-28-0	Thallium	0.78	U		F
7440-62-2	Vanadium	23.30			P
7440-66-6	Zinc	384.00			P
	Cyanide	0.98	U		AS

Color Before: BROWN

Clarity Before:

Texture: COARSE

Color After: LT.BROWN

Clarity After:

Artifacts:

Comments:

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE12

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302003

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 50.4

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	3190.00			P
7440-36-0	Antimony	11.90	U		P
7440-38-2	Arsenic	6.10			F
7440-39-3	Barium	118.00			P
7440-41-7	Beryllium	1.30	B		P
7440-43-9	Cadmium	1.40	B		P
7440-70-2	Calcium	4070.00			P
7440-47-3	Chromium	31.50			P
7440-48-4	Cobalt	32.90			P
7440-50-8	Copper	58.30			P
7439-89-6	Iron	17800.00			P
7439-92-1	Lead	6.30		S	F
7439-95-4	Magnesium	3110.00			P
7439-96-5	Manganese	1230.00			P
7439-97-6	Mercury	0.30			CV
7439-02-0	Nickel	40.40			P
7440-09-7	Potassium	560.00	B		P
7782-49-2	Selenium	0.79	U	+	F
7440-22-4	Silver	0.79	U		P
7440-23-5	Sodium	221.00	B		P
7440-28-0	Thallium	0.79	U		F
7440-62-2	Vanadium	22.80			P
7440-66-6	Zinc	288.00			P
	Cyanide	0.99	U		AS

Color Before: BROWN

Clarity Before:

Texture: COARSE

Color After: LT.BROWN

Clarity After:

Artifacts:

Comments:

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE13

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302004

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 60.6

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum_	12400.00			P
7440-36-0	Antimony_	9.90	U		P
7440-38-2	Arsenic__	7.80			F
7440-39-3	Barium__	146.00			P
7440-41-7	Beryllium	2.20			P
7440-43-9	Cadmium__	0.99	U		P
7440-70-2	Calcium__	1800.00			P
7440-47-3	Chromium_	19.30			P
7440-48-4	Cobalt__	31.00			P
7440-50-8	Copper__	61.00			P
7439-89-6	Iron_____	17500.00			P
7439-92-1	Lead_____	3.30		S	F
7439-95-4	Magnesium	2580.00			P
7439-96-5	Manganese	291.00			P
7439-97-6	Mercury__	0.50			CV
7439-02-0	Nickel__	55.20			P
7440-09-7	Potassium	869.00	B		P
7782-49-2	Selenium_	0.96	B	W	F
7440-22-4	Silver__	0.66	U		P
7440-23-5	Sodium__	80.60	B		P
7440-28-0	Thallium_	0.66	U		F
7440-62-2	Vanadium_	23.20			P
7440-66-6	Zinc_____	299.00			P
	Cyanide__	0.83	U		AS

Color Before: BROWN

Clarity Before:

Texture: COARSE

Color After: LT.BROWN

Clarity After:

Artifacts:

Comments:

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE14

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302005

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 43.2

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	10500.00			P
7440-36-0	Antimony	13.90	U		P
7440-38-2	Arsenic	3.80	B	S	F
7440-39-3	Barium	154.00			P
7440-41-7	Beryllium	2.10	B		P
7440-43-9	Cadmium	2.20	B		P
7440-70-2	Calcium	4560.00			P
7440-47-3	Chromium	43.20			P
7440-48-4	Cobalt	59.70			P
7440-50-8	Copper	105.00			P
7439-89-6	Iron	23000.00			P
7439-92-1	Lead	5.30		S	F
7439-95-4	Magnesium	3100.00			P
7439-96-5	Manganese	1540.00			P
7439-97-6	Mercury	0.35			CV
7439-02-0	Nickel	75.20			P
7440-09-7	Potassium	970.00	B		P
7782-49-2	Selenium	0.93	U	S	F
7440-22-4	Silver	0.93	U		P
7440-23-5	Sodium	195.00	B		P
7440-28-0	Thallium	0.93	U		F
7440-62-2	Vanadium	28.20			P
7440-66-6	Zinc	497.00			P
	Cyanide	1.20	U		AS

Color Before: DK.BROWN

Clarity Before:

Texture: COARSE

Color After: LT.YELLOW

Clarity After:

Artifacts:

Comments:

U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE15

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302006

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 83.6

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum_	6810.00			P
7440-36-0	Antimony_	7.20	U		P
7440-38-2	Arsenic_	372.00		S	P
7440-39-3	Barium_	158.00			P
7440-41-7	Beryllium_	0.70	B		P
7440-43-9	Cadmium_	2.60			P
7440-70-2	Calcium_	1270.00			P
7440-47-3	Chromium_	29.10			P
7440-48-4	Cobalt_	14.30			P
7440-50-8	Copper_	42.70			P
7439-89-6	Iron_	37700.00			P
7439-92-1	Lead_	3.50		S	P
7439-95-4	Magnesium_	1830.00			P
7439-96-5	Manganese_	411.00			P
7439-97-6	Mercury_	0.18			CV
7439-02-0	Nickel_	20.90			P
7440-09-7	Potassium_	322.00	B		P
7782-49-2	Selenium_	2.40	U	S	P
7440-22-4	Silver_	0.48	U		P
7440-23-5	Sodium_	129.00	B		P
7440-28-0	Thallium_	0.48	U		P
7440-62-2	Vanadium_	41.10			P
7440-66-6	Zinc_	1230.00			P
	Cyanide_	0.68			AS

Color Before: BROWN

Clarity Before:

Texture: COARSE

Color After: LT.YELLOW

Clarity After:

Artifacts:

Comments:

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE16

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302007

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 43.7

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	2140.00			P
7440-36-0	Antimony	13.70	U		P
7440-38-2	Arsenic	62.40		+	F
7440-39-3	Barium	407.00			P
7440-41-7	Beryllium	0.55	B		P
7440-43-9	Cadmium	2.60			P
7440-70-2	Calcium	2220.00	B		P
7440-47-3	Chromium	63.90			P
7440-48-4	Cobalt	9.60	B		P
7440-50-8	Copper	71.50			P
7439-89-6	Iron	39000.00			P
7439-92-1	Lead	4810.00			F
7439-95-4	Magnesium	784.00	B		P
7439-96-5	Manganese	464.00			P
7439-97-6	Mercury	2.90			CV
7439-02-0	Nickel	10.60	B		P
7440-09-7	Potassium	330.00	U		P
7782-49-2	Selenium	0.92	U		F
7440-22-4	Silver	0.92	U		P
7440-23-5	Sodium	52.40	B		P
7440-28-0	Thallium	0.92	U		F
7440-62-2	Vanadium	26.50			P
7440-66-6	Zinc	651.00			P
	Cyanide	1.80			AS

Color Before: BLACK

Clarity Before:

Texture: COARSE

Color After: LT.YELLOW

Clarity After:

Artifacts:

Comments:



## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE17

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302008

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 81.8

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	4970.00			P
7440-36-0	Antimony	7.30	U		P
7440-38-2	Arsenic	3.90		S	P
7440-39-3	Barium	80.60			P
7440-41-7	Beryllium	0.71	B		P
7440-43-9	Cadmium	0.73	U		P
7440-70-2	Calcium	1520.00			P
7440-47-3	Chromium	12.40			P
7440-48-4	Cobalt	8.60	B		P
7440-50-8	Copper	15.40			P
7439-89-6	Iron	12800.00			P
7439-92-1	Lead	1.90		S	P
7439-95-4	Magnesium	1040.00	B		P
7439-96-5	Manganese	811.00			P
7439-97-6	Mercury	0.12	U		CV
7439-02-0	Nickel	9.80	B		P
7440-09-7	Potassium	782.00	B		P
7782-49-2	Selenium	0.71	B	W	P
7440-22-4	Silver	0.49	U		P
7440-23-5	Sodium	64.50	B		P
7440-28-0	Thallium	0.49	U		P
7440-62-2	Vanadium	15.70			P
7440-66-6	Zinc	200.00			P
	Cyanide	0.61	U		AS

Color Before: RED-BROWN

Clarity Before:

Texture: COARSE

Color After: LT.BROWN

Clarity After:

Artifacts:

Comments:

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE18

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302009

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 76.0

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	3400.00			P
7440-36-0	Antimony	7.90	U		P
7440-38-2	Arsenic	6.70			F
7440-39-3	Barium	52.70			P
7440-41-7	Beryllium	0.51	B		P
7440-43-9	Cadmium	0.79	U		P
7440-70-2	Calcium	2110.00			P
7440-47-3	Chromium	10.60			P
7440-48-4	Cobalt	6.40	B		P
7440-50-8	Copper	10.00			P
7439-89-6	Iron	10800.00			P
7439-92-1	Lead	182.00			F
7439-95-4	Magnesium	937.00	B		P
7439-96-5	Manganese	318.00			P
7439-97-6	Mercury	0.13	U		CV
7439-02-0	Nickel	7.40	B		P
7440-09-7	Potassium	397.00	B		P
7782-49-2	Selenium	0.63	B	W	F
7440-22-4	Silver	0.53	U		P
7440-23-5	Sodium	43.50	B		P
7440-28-0	Thallium	0.53	U		F
7440-62-2	Vanadium	19.60			P
7440-66-6	Zinc	169.00			P
	Cyanide	0.66	U		AS

Color Before: RED-BROWN

Clarity Before:

Texture: COARSE

Color After: LT.BROWN

Clarity After:

Artifacts:

Comments:

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE19

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302012

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 83.1

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	4170.00			P
7440-38-0	Antimony	7.20	U		P
7440-38-2	Arsenic	5.40			F
7440-39-3	Barium	66.00			P
7440-41-7	Beryllium	0.58	B		P
7440-43-9	Cadmium	0.72	U		P
7440-70-2	Calcium	1620.00			P
7440-47-3	Chromium	8.10			P
7440-48-4	Cobalt	7.50	B		P
7440-50-8	Copper	7.10			P
7439-89-6	Iron	9880.00			P
7439-92-1	Lead	16.90			F
7439-95-4	Magnesium	1140.00	B		P
7439-96-5	Manganese	351.00			P
7439-97-6	Mercury	0.12	U		CV
7439-02-0	Nickel	7.10	B		P
7440-09-7	Potassium	483.00	B		P
7782-49-2	Selenium	0.60	B		F
7440-22-4	Silver	0.48	U		P
7440-23-5	Sodium	115.00	B		P
7440-28-0	Thallium	0.48	U		F
7440-62-2	Vanadium	12.30			P
7440-66-6	Zinc	201.00			P
	Cyanide	0.60	U		AS

Color Before: RED-BROWN

Clarity Before:

Texture: COARSE

Color After: LT.BROWN

Clarity After:

Artifacts:

Comments:

10/15/92  
Rd.

U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE20

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302013

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 64.7

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum_	3760.00			P
7440-38-0	Antimony_	9.30	U		P
7440-38-2	Arsenic_	9.00			P
7440-39-3	Barium_	80.90			P
7440-41-7	Beryllium	0.53	B		P
7440-43-9	Cadmium_	1.10	B		P
7440-70-2	Calcium_	12400.00			P
7440-47-3	Chromium_	38.60			P
7440-48-4	Cobalt_	8.80	B		P
7440-50-8	Copper_	81.70			P
7439-89-6	Iron_	10300.00			P
7439-92-1	Lead_	253.00			P
7439-95-4	Magnesium	4750.00			P
7439-96-5	Manganese	437.00			P
7439-97-6	Mercury_	0.23			CV
7439-02-0	Nickel_	10.70	B		P
7440-09-7	Potassium	534.00	B		P
7782-49-2	Selenium_	0.96	B	W	P
7440-22-4	Silver_	0.62	U		P
7440-23-5	Sodium_	63.40	B		P
7440-28-0	Thallium_	0.62	U		P
7440-62-2	Vanadium_	12.70	B		P
7440-66-6	Zinc_	226.00			P
	Cyanide_	0.77	U		AS

Color Before: BROWN

Clarity Before:

Texture: COARSE

Color After: LT.BROWN

Clarity After:

Artifacts:

Comments:

## U.S. EPA - CLP

1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE21

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302014

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 74.9

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	5090.00			P
7440-36-0	Antimony	11.00	B		P
7440-38-2	Arsenic	33.30		S	F
7440-39-3	Barium	268.00			P
7440-41-7	Beryllium	0.65	B		P
7440-43-9	Cadmium	2.30			P
7440-70-2	Calcium	17600.00			P
7440-47-3	Chromium	45.10			P
7440-48-4	Cobalt	14.80			P
7440-50-8	Copper	388.00			P
7439-89-6	Iron	67800.00			P
7439-92-1	Lead	358.00			F
7439-95-4	Magnesium	4860.00			P
7439-96-5	Manganese	801.00			P
7439-97-6	Mercury	0.33			CV
7439-02-0	Nickel	56.50			P
7440-09-7	Potassium	1100.00	B		P
7782-49-2	Selenium	0.53	U	W	F
7440-22-4	Silver	0.53	U		P
7440-23-5	Sodium	219.00	B		P
7440-28-0	Thallium	0.53	U		F
7440-62-2	Vanadium	73.10			P
7440-66-6	Zinc	1150.00			P
	Cyanide	0.67	U		AS

Color Before: BROWN

Clarity Before:

Texture: COARSE

Color After: LT.YELLOW

Clarity After:

Artifacts:

Comments:

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1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE22

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302015

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 74.8

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	5470.00			P
7440-36-0	Antimony	8.00	U		P
7440-38-2	Arsenic	10.40			P
7440-39-3	Barium	113.00			P
7440-41-7	Beryllium	0.78	B		P
7440-43-9	Cadmium	0.80	U		P
7440-70-2	Calcium	2670.00			P
7440-47-3	Chromium	8.20			P
7440-48-4	Cobalt	8.70	B		P
7440-50-8	Copper	17.00			P
7439-89-6	Iron	11600.00			P
7439-92-1	Lead	22.20		W	P
7439-95-4	Magnesium	1070.00	B		P
7439-96-5	Manganese	469.00			P
7439-97-6	Mercury	0.13	U		CV
7439-02-0	Nickel	12.70			P
7440-09-7	Potassium	483.00	B		P
7782-49-2	Selenium	0.72	B	W	P
7440-22-4	Silver	0.53	U		P
7440-23-5	Sodium	91.00	B		P
7440-28-0	Thallium	0.53	U		P
7440-62-2	Vanadium	19.80			P
7440-66-6	Zinc	286.00			P
	Cyanide	0.67	U		AS

Color Before: RED-BROWN

Clarity Before:

Texture: COARSE

Color After: LT.BROWN

Clarity After:

Artifacts:

Comments:

ORIGINAL  
File

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1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE23

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302016

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 77.4

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum_	8230.00			P
7440-36-0	Antimony_	7.80	U		P
7440-38-2	Arsenic_	5.60			P
7440-39-3	Barium_	111.00			P
7440-41-7	Beryllium_	1.00	B		P
7440-43-9	Cadmium_	0.78	U		P
7440-70-2	Calcium_	1730.00			P
7440-47-3	Chromium_	15.20			P
7440-48-4	Cobalt_	11.10	B		P
7440-50-8	Copper_	20.10			P
7439-89-6	Iron_	20000.00			P
7439-92-1	Lead_	20.20			P
7439-95-4	Magnesium_	1890.00			P
7439-96-5	Manganese_	1090.00			P
7439-97-6	Mercury_	0.13	U		CV
7439-02-0	Nickel_	14.40			P
7440-09-7	Potassium_	383.00	B		P
7782-49-2	Selenium_	0.52	U		P
7440-22-4	Silver_	0.52	U		P
7440-23-5	Sodium_	41.10	B		P
7440-28-0	Thallium_	0.52	U		P
7440-62-2	Vanadium_	31.50			P
7440-66-6	Zinc_	61.80			P
	Cyanide_	0.65	U		AS

Color Before: BROWN

Clarity Before:

Texture: COARSE

Color After: LT.YELLOW

Clarity After:

Artifacts:

Comments:

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## INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE24

b Name: KEYSTONE LAB-HOUSTON

Contract: 68-D0-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE02

Matrix (soil/water): WATER

Lab Sample ID: 9201301013

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): ug/L

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	15.10	B	E	P
7440-36-0	Antimony	30.00	U		P
7440-38-2	Arsenic	5.40	B		F
7440-39-3	Barium	1.00	U		P
7440-41-7	Beryllium	1.00	U		P
7440-43-9	Cadmium	3.00	U		P
7440-70-2	Calcium	29.10	B		P
7440-47-3	Chromium	3.00	U		P
7440-48-4	Cobalt	4.00	U		P
7440-50-8	Copper	3.00	U		P
7439-89-6	Iron	12.50	B	NE	P
7439-92-1	Lead	1.00	U		F
7439-95-4	Magnesium	42.60	B		P
7439-96-5	Manganese	2.00	U		P
7439-97-6	Mercury	0.20	U		CV
7439-02-0	Nickel	22.00	U		P
7440-09-7	Potassium	722.00	U		P
7782-49-2	Selenium	2.00	U		F
7440-22-4	Silver	2.00	U		P
7440-23-5	Sodium	30.00	U		P
7440-28-0	Thallium	2.00	U		F
7440-62-2	Vanadium	4.00	U		P
7440-66-6	Zinc	15.80	B		P
	Cyanide	10.00	U		AS

Color Before: COLORLESS

Clarity Before: CLEAR

Texture:

Color After: COLORLESS

Clarity After: CLEAR

Artifacts:

Comments:



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1  
INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MCJE26

Lab Name: KEYSTONE LAB-HOUSTON

Contract: 68-DO-0147

Lab Code: KEYTX

Case No.: 17744

SAS No.:

SDG No.: MCJE11

Matrix (soil/water): SOIL

Lab Sample ID: 9201302017

Level (low/med): LOW

Date Received: 01/29/92

% Solids: 67.9

Concentration Units (ug/L or mg/kg dry weight): MG/KG

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum_	6790.00			P
7440-36-0	Antimony_	8.80	U		P
7440-38-2	Arsenic_	3.40		W	P
7440-39-3	Barium_	44.60	B		P
7440-41-7	Beryllium	0.72	B		P
7440-43-9	Cadmium_	0.88	U		P
7440-70-2	Calcium_	382.00	B		P
7440-47-3	Chromium_	14.20			P
7440-48-4	Cobalt_	9.80	B		P
7440-50-8	Copper_	13.50			P
7439-89-6	Iron_	18100.00			P
7439-92-1	Lead_	17.80			P
7439-95-4	Magnesium	1250.00	B		P
7439-96-5	Manganese	366.00			P
7439-97-6	Mercury_	0.15	U		CV
7439-02-0	Nickel_	10.70	B		P
7440-09-7	Potassium	582.00	B		P
7782-49-2	Selenium_	0.82	B	M	P
7440-22-4	Silver_	0.59	U		P
7440-23-5	Sodium_	32.10	B		P
7440-28-0	Thallium_	0.59	U		P
7440-62-2	Vanadium_	27.20			P
7440-66-6	Zinc_	129.00			P
	Cyanide_	0.74	U		AS

Color Before: BROWN

Clarity Before:

Texture: COARSE

Color After: LT.YELLOW

Clarity After:

Artifacts:

Comments:

## **APPENDIX C**

## HOME WELL SURVEY

Home Owner's Name: (b) (6)

Date: \_\_\_\_\_

Address: (b) (6)

Home Phone: (b) (6)

Work Phone: \_\_\_\_\_

- 1 Please describe the type of home well you presently utilize:  
(Check those which apply)

\_\_\_\_\_ Dug well  
☒ Drilled by a rig; if so, please identify company (name, address, and phone):  
Gardner - Sassafrasville  
\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_ Other (describe) \_\_\_\_\_

- 1a. Please estimate the following:

Year installed unknownDate of last service 6/15/92

Company who serviced (name, address, and phone):

New Pump  
\_\_\_\_\_  
\_\_\_\_\_

2. Please provide the following measurements of your well:

a.

b.

(b) (9)

3. Please describe the casing material used in your well:

a. Composition

☒ Iron☐ PVC☐ Galvanized☐ Terra Cotta☐ Other - Please

Specify (if known)

b. Length (if known): unknown

GRU 100  
10000

## HOME WELL SURVEY

Home Owner's Name: (b) (6) Date: \_\_\_\_\_

4 Please describe, if known, any screening material used in your well:

a. Length of screen: \_\_\_\_\_

b. Depth of screen in well: \_\_\_\_\_

5 Please indicate, if known, the depth to the groundwater in your well (from the surface):

unknown

6 Please indicate the composition of home plumbing (pipes) in your system:

\_\_\_\_\_ Iron \_\_\_\_\_ PVC \_\_\_\_\_ Galvanized \_\_\_\_\_ Lead

✓ Other (describe): Copper

7 Please describe the water pump used in your system:

a. Location of the pump

✓ Inside the well (submersible pump); Depth in well: \_\_\_\_\_

\_\_\_\_\_ Outside the well (indicate location): \_\_\_\_\_

b. Type of pump

Branch (if known): \_\_\_\_\_

Capacity (gallons per minute): \_\_\_\_\_

c. Estimate hours of pump operation per day: \_\_\_\_\_

d. Is storage tank used: ✓ Yes \_\_\_\_\_ No

Type (material) bladder Capacity \_\_\_\_\_

8 a. Do you regularly or have you ever added chemicals directly to your well?  
(i.e., chlorine, clorox, etc.) \_\_\_\_\_ Yes ✓ No

If yes, date last added: \_\_\_\_\_ Approximate amount added \_\_\_\_\_

Compound (brand name): \_\_\_\_\_

## HOME WELL SURVEY

Home Owner's Name: \_\_\_\_\_

(b) (6)

Date: \_\_\_\_\_

- b. Please describe any type of water treatment you are currently using (check those which apply):

\_\_\_\_\_ Filtration

\_\_\_\_\_ Other (explain)

Type: \_\_\_\_\_

\_\_\_\_\_ none

\_\_\_\_\_ Water Softeners

Indicate Brand: \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

9. Please indicate any testing that has been done on your water:

Date of testing: none

Name of individual(s) responsible for testing: \_\_\_\_\_

10. Well Use: ✓ Drinking

✓ Other: household

11. Do you notice color, taste, or odor problems with well water? \_\_\_\_\_ Yes ✓ No

If yes, identify: \_\_\_\_\_

Do you notice water supply problems? \_\_\_\_\_ Yes

✓ No

If yes, when: \_\_\_\_\_ how often: \_\_\_\_\_

12. Please indicate the type(s) of wastewater system used (check):

Sewer Line ✓

Septic Tank \_\_\_\_\_

Cesspool \_\_\_\_\_

Drain Field \_\_\_\_\_

Distance to Well \_\_\_\_\_

13. We may be taking water samples from many area homes in the near future. If your well is chosen for sampling, would you be willing to allow our NUS representatives to sample your well? Sampling involves collecting water from one of your indoor or outdoor spigots.

✓ Yes, I will allow my well to be sampled.

\_\_\_\_\_ No, I will not allow my well to be sampled.

## HOME WELL SURVEY

Home Owner's Name: (b) (6) Date: \_\_\_\_\_

If yes, please indicate the time of day which would be convenient for us to sample.

\_\_\_\_\_ Morning \_\_\_\_\_ Afternoon \_\_\_\_\_ Evening

14. In the space below, please furnish a rough sketch of your property, indicating the location of your well and on-lot wastewater system, if applicable. Also indicate the location of the spigot you would prefer us to sample.

outside faucet.

HW-2

# HOME WELL SURVEY

Home Owner's Name: (b) (6)

Date: 1/28/92

Address: (b) (6)

Home Phone: (b) (6)

Work Phone:

1. Please describe the type of home well you presently utilize:  
(Check those which apply)

☐ Dug well

☒ Drilled by a rig; if so, please identify company (name, address, and phone):

\_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

☐ Other (describe) \_\_\_\_\_

- 1a. Please estimate the following: Year installed \_\_\_\_\_

Date of last service 65

Company who serviced (name, address, and phone): \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

2. Please provide the following measurements of your well:

a. (b) (9)  
b. (b) (9)

3. Please describe the casing material used in your well:

a. Composition

☐ Iron

☒ PVC

☐ Galvanized

☐ Terra Cotta

☐ Other - Please

Specify (if known)

b. Length (if known): \_\_\_\_\_

## HOME WELL SURVEY

Home Owner's Name: \_\_\_\_\_

(b) (6)

Date: \_\_\_\_\_

4. Please describe, if known, any screening material used in your well:

a. Length of screen: \_\_\_\_\_

b. Depth of screen in well: \_\_\_\_\_

5. Please indicate, if known, the depth to the groundwater in your well (from the surface):

55

6. Please indicate the composition of home plumbing (pipes) in your system:

☒ Iron    ☐ PVC    ☐ Galvanized    ☐ Lead  
☐ Other (describe): \_\_\_\_\_

7. Please describe the water pump used in your system:

a. Location of the pump

☒ Inside the well (submersible pump); Depth in well: \_\_\_\_\_

\_\_\_\_\_ Outside the well (indicate location): \_\_\_\_\_

b. Type of pump

Branch (if known): \_\_\_\_\_

Capacity (gallons per minute): \_\_\_\_\_

c. Estimate hours of pump operation per day: \_\_\_\_\_

d. Is storage tank used: ☒ Yes    ☐ NoType (material) \_\_\_\_\_ Capacity 50

8. a. Do you regularly or have you ever added chemicals directly to your well?

(i.e., chlorine, clorox, etc.) ☐ Yes    ☒ No

If yes, date last added: \_\_\_\_\_ Approximate amount added \_\_\_\_\_

Compound (brand name): \_\_\_\_\_



# HOME WELL SURVEY

Home Owner's Name: \_\_\_\_\_

**(b) (6)**

Date: \_\_\_\_\_

- b. Please describe any type of water treatment you are currently using (check those which apply):

\_\_\_\_\_ Filtration

\_\_\_\_\_ Other (explain)

Type: \_\_\_\_\_

\_\_\_\_\_ Water Softeners

Indicate Brand: \_\_\_\_\_

9. Please indicate any testing that has been done on your water:

Date of testing: \_\_\_\_\_

Name of individual(s) responsible for testing: \_\_\_\_\_

10. Well Use: ☒ Drinking \_\_\_\_\_ Other: \_\_\_\_\_

11. Do you notice color, taste, or odor problems with well water? \_\_\_\_\_ Yes ☒ No

If yes, identify: \_\_\_\_\_

Do you notice water supply problems? \_\_\_\_\_ Yes ☒ No

If yes, when: \_\_\_\_\_ how often: \_\_\_\_\_

12. Please indicate the type(s) of wastewater system used (check):

Sewer Line ☒ \_\_\_\_\_

Septic Tank \_\_\_\_\_

Cesspool \_\_\_\_\_

Drain Field \_\_\_\_\_

Distance to Well \_\_\_\_\_

13. We may be taking water samples from many area homes in the near future. If your well is chosen for sampling, would you be willing to allow our NUS representatives to sample your well? Sampling involves collecting water from one of your indoor or outdoor spigots.

\_\_\_\_\_ Yes, I will allow my well to be sampled.

\_\_\_\_\_ No, I will not allow my well to be sampled.

# HOME WELL SURVEY

Home Owner's Name: (b) (6) \_\_\_\_\_

Date: \_\_\_\_\_

If yes, please indicate the time of day which would be convenient for us to sample.

\_\_\_\_\_ Morning \_\_\_\_\_ Afternoon \_\_\_\_\_ Evening

14. In the space below, please furnish a rough sketch of your property, indicating the location of your well and on-lot wastewater system, if applicable. Also indicate the location of the spigot you would prefer us to sample.

**PLATE 1**

EPA REGION III  
SUPERFUND DOCUMENT MANAGEMENT SYSTEM

DOC ID # 404788  
PAGE #           

**IMAGERY COVER SHEET**  
**UNSCANNABLE ITEM**

Contact the CERCLA Records Center to view this document.

SITE NAME	<u>Linfield Industrial</u>
OPERABLE UNIT	<u>00</u>
SECTION/BOX/FOLDER	<u>1C - 1 - 1.002</u>

REPORT OR DOCUMENT TITLE	<u>Final Screening</u> <u>Site Inspection</u>
DATE OF DOCUMENT	<u>Sept. 1, 1992</u>
DESCRIPTION OF IMAGERY	<u>Four Mile Radius</u> <u>MAP</u>
NUMBER AND TYPE OF IMAGERY ITEM(S)	<u>1 oversized map</u>